

(QT Reviewed)

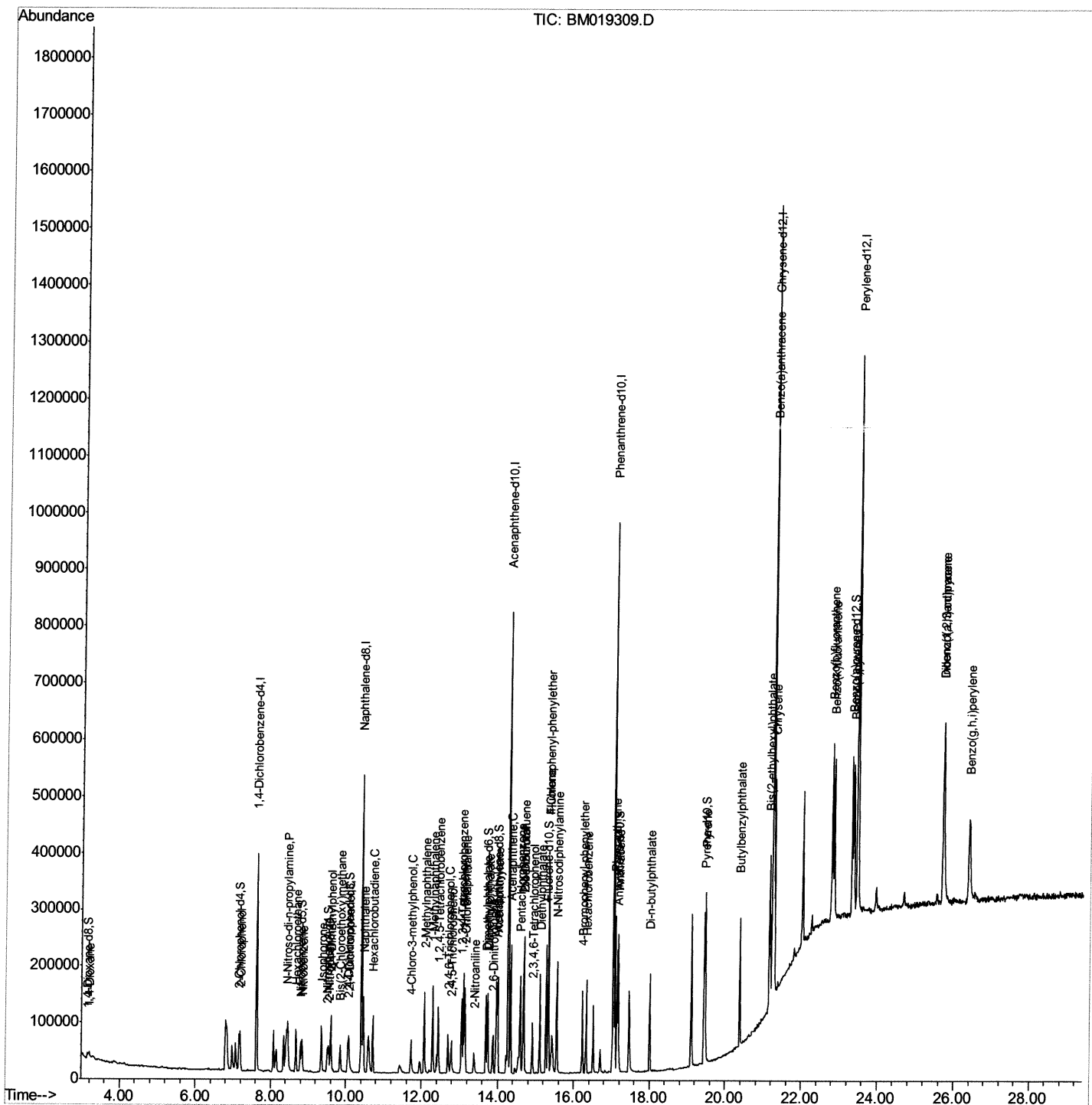
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Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\  
Data File : BM019309.D  
Acq On    : 22 Mar 2019 12:49  
Operator  : JU/SJ  
Sample    : SSTD00552  
Misc      :  
ALS Vial  : 2      Sample Multiplier: 1
```

Instrument :
BNA_M
ClientSampleId :
SSTD00552

Manual Integrations

Sohil
3/25/2019 6:35:51 PM

Quant Time: Mar 22 16:24:07 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 22 16:14:04 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

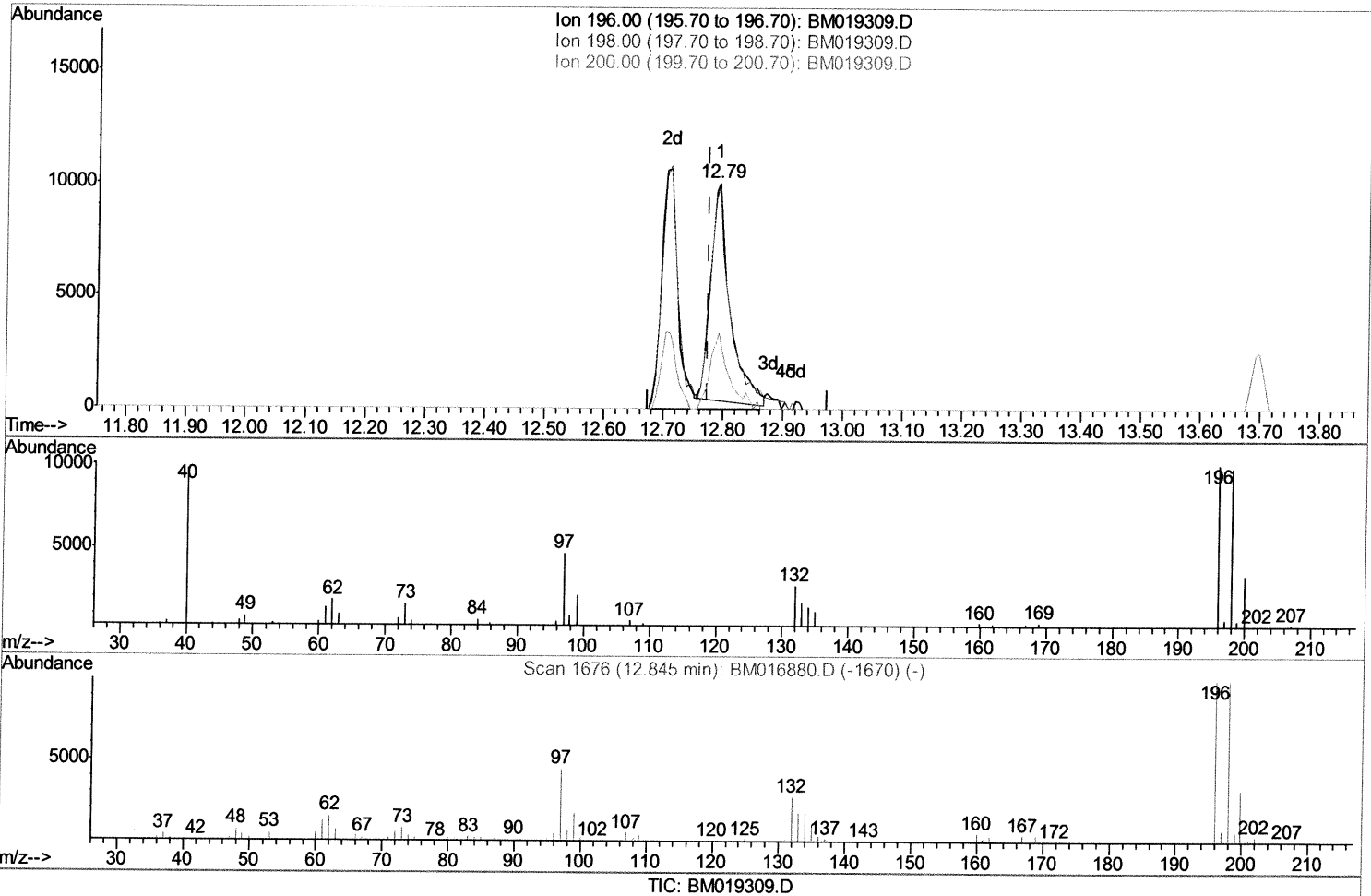
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
 Data File : BM019309.D
 Acq On : 22 Mar 2019 12:49
 Operator : JU/SJ
 Sample : SST00552
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST00552

Manual Integrations
 APPROVED

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 3/25/2019 6:35:51 PM

Quant Time: Mar 22 16:22:17 2019
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 22 16:14:04 2019
 Response via : Initial Calibration



(40) 2,4,5-Trichlorophenol
 12.792min (+0.018) 4.03ng/ul
 response 21637

Ion	Exp%	Act%
196.00	100	100
198.00	93.40	97.80
200.00	31.70	33.91
0.00	0.00	0.00

Quantitation Report (Qedit)

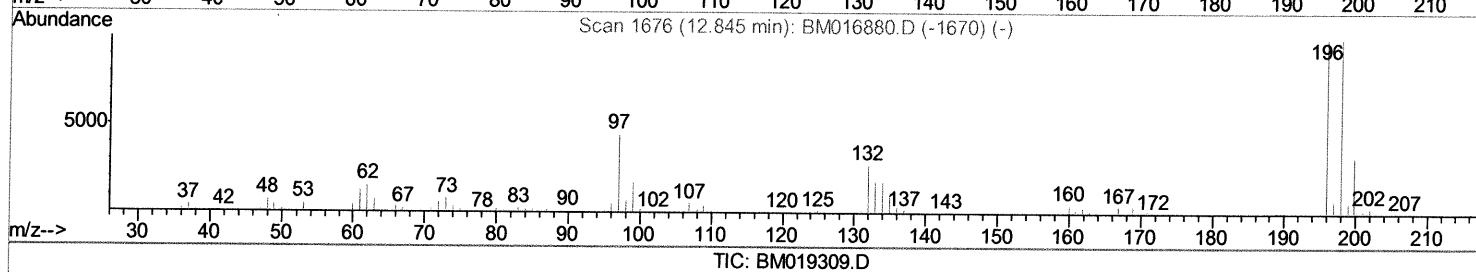
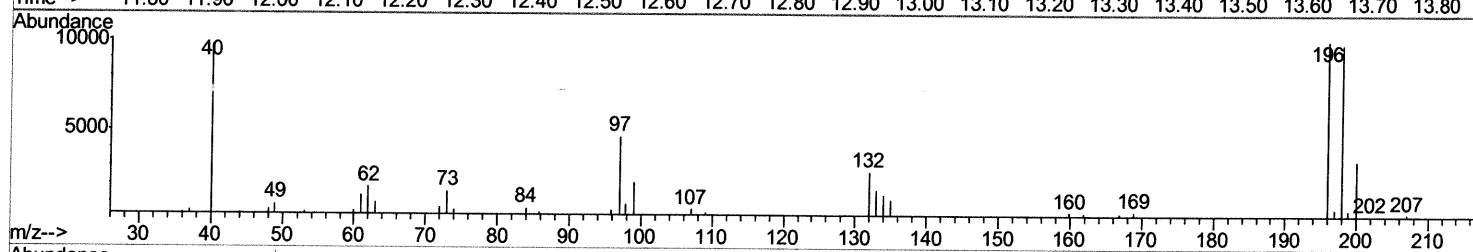
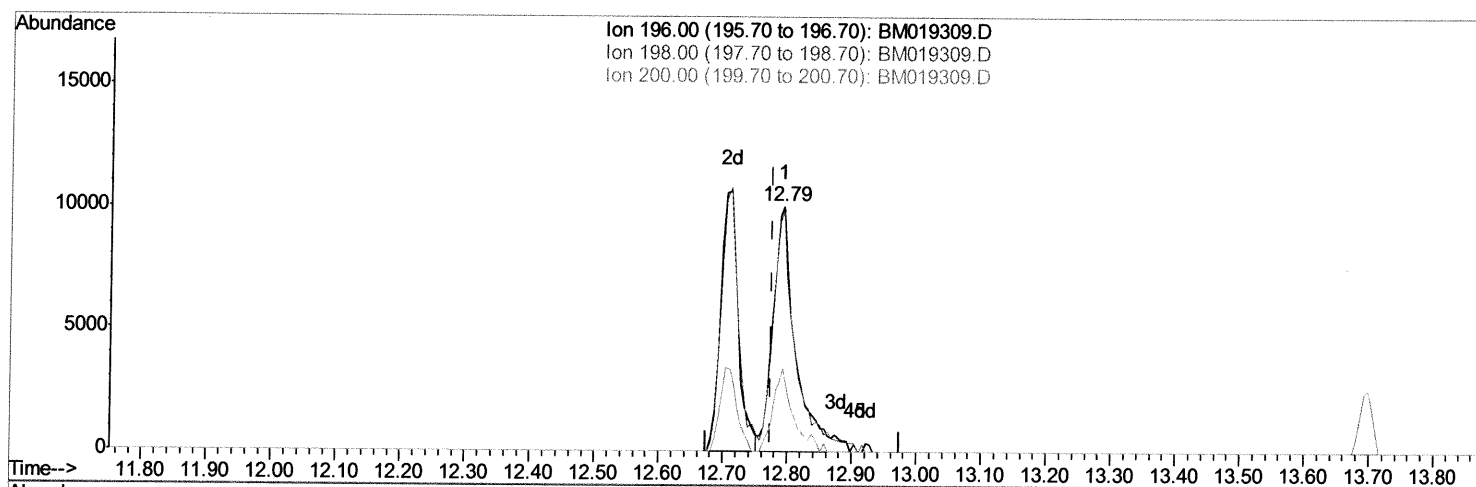
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(40) 2,4,5-Trichlorophenol

12.792min (+0.018) 4.58ng/ul m

> 51 3/29/19

response 24620

Ion	Exp%	Act%
196.00	100	100
198.00	93.40	97.80
200.00	31.70	33.91
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	105016	20.00	nq/ul	0.00
18) Naphthalene-d8	10.40	136	431783	20.00	nq/ul	0.00
36) Acenaphthene-d10	14.27	164	251668	20.00	nq/ul	0.00
62) Phenanthrene-d10	17.03	188	559984	20.00	nq/ul	0.00
78) Chrysene-d12	21.24	240	638027	20.00	nq/ul	0.00
86) Perylene-d12	23.46	264	700830	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	5282	2.15	nq/uL	0.00
5) Phenol-d5	0.00	99	0d	0.00	nq/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	nq/ul	
9) 2-Chlorophenol-d4	7.15	132	31493	4.36	nq/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	nq/ul	
19) Nitrobenzene-d5	8.79	128	13219	4.66	nq/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	13590	5.27	nq/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	27635	4.22	nq/ul	0.01
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	nq/ul	
44) Dimethylphthalate-d6	13.70	166	97458	4.82	nq/ul	0.00
47) Acenaphthylene-d8	13.97	160	115570	4.43	nq/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	nq/ul	
58) Fluorene-d10	15.27	176	87725	4.97	nq/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	nq/ul	
71) Anthracene-d10	17.13	188	129171	4.78	nq/ul	0.00
79) Pyrene-d10	19.44	212	139918	4.52	nq/ul	0.00
90) Benzo(a)pyrene-d12	23.32	264	172653	4.72	ng/ul	0.00

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.19	88	6150	2.407	nq/uL	95
10) 2-Chlorophenol	7.19	128	31839	4.385	nq/ul	95
15) N-Nitroso-di-n-propylamine	8.43	70	27816	5.233	nq/ul#	85
17) Hexachloroethane	8.67	117	14257	4.999	nq/ul	84
20) Nitrobenzene	8.83	77	42753	5.957	nq/ul	98
21) Isophorone	9.35	82	71398	5.006	nq/ul	95
23) 2-Nitrophenol	9.54	139	16286	5.469	nq/ul	95
24) 2,4-Dimethylphenol	9.59	107	39747	5.173	nq/ul	96
25) Bis(2-Chloroethoxy)methane	9.84	93	43027	4.881	nq/ul	98
27) 2,4-Dichlorophenol	10.07	162	26142	4.151	nq/ul	94
28) Naphthalene	10.45	128	110927	4.996	nq/ul	98
31) Hexachlorobutadiene	10.71	225	20701	4.870	nq/ul	98
33) 4-Chloro-3-methylphenol	11.72	107	30316	4.593	nq/ul	99
34) 2-Methylnaphthalene	12.07	142	77250	4.938	nq/ul	100
35) 1-Methylnaphthalene	12.30	142	74758	4.779	nq/ul#	98
37) 1,2,4,5-Tetrachlorobenzene	12.45	216	39032	4.516	nq/ul#	97
39) 2,4,6-Trichlorophenol	12.71	196	20117	4.087	nq/ul	95
40) 2,4,5-Trichlorophenol	12.79	196	24620m)	4.584	nq/ul	
41) 1,1'-Biphenyl	13.10	154	101892	4.867	nq/ul#	94
42) 2-Chloronaphthalene	13.15	162	76622	4.676	nq/ul	97
43) 2-Nitroaniline	13.39	65	16411	4.547	nq/ul	89
45) Dimethylphthalate	13.75	163	100615	5.095	nq/ul#	96
46) 2,6-Dinitrotoluene	13.88	165	15135	4.990	ng/ul	93

SJ 3/29/19

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48) Acenaphthylene	14.00	152	116972	4.735	nq/ul	98
50) Acenaphthene	14.34	153	88071	5.021	nq/ul	99
54) Dibenzofuran	14.68	168	121488	5.049	nq/ul	98
55) 2,4-Dinitrotoluene	14.67	165	21526	4.826	nq/ul#	34
56) 2,3,4,6-Tetrachlorophenol	14.92	232	19353	4.455	nq/ul#	82
57) Diethylphthalate	15.11	149	96647	4.940	nq/ul	96
59) Fluorene	15.33	166	96429	4.878	nq/ul	99
60) 4-Chlorophenyl-phenylether	15.33	204	49313	4.920	nq/ul	92
65) N-Nitrosodiphenylamine	15.55	169	81381	4.698	nq/ul	98
66) 4-Bromophenyl-phenylether	16.23	248	28765	4.540	nq/ul	95
67) Hexachlorobenzene	16.33	284	37243	5.462	nq/ul#	89
70) Phenanthrene	17.07	178	157683	5.069	nq/ul	99
72) Anthracene	17.17	178	154900	4.851	nq/ul	99
73) 1,2,3,4-Tetrachlorobenzene	13.06	216	38072	4.108	nq/uL	98
74) Pentachlorobenzene	14.59	250	41173	4.925	nq/uL	98
76) Di-n-butylphthalate	18.00	149	135340	4.144	nq/ul	98
80) Pyrene	19.47	202	183093	4.674	nq/ul#	87
81) Butylbenzylphthalate	20.37	149	60890	4.203	nq/ul#	86
83) Benzo(a)anthracene	21.22	228	184310	4.666	nq/ul	100
84) Bis(2-ethylhexyl)phthalate	21.14	149	85578	3.758	nq/ul	97
85) Chrysene	21.27	228	184079	4.997	nq/ul	99
88) Benzo(b)fluoranthene	22.79	252	198292	4.727	nq/ul#	96
89) Benzo(k)fluoranthene	22.84	252	197970	4.915	nq/ul#	96
91) Benzo(a)pyrene	23.36	252	188167	4.706	nq/ul#	96
92) Indeno(1,2,3-cd)pyrene	25.70	276	235314	4.933	nq/ul#	90
93) Dibenzo(a,h)anthracene	25.71	278	198861	4.984	nq/ul#	95
94) Benzo(a,h,i)perylene	26.39	276	200273	5.064	nq/ul#	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed