

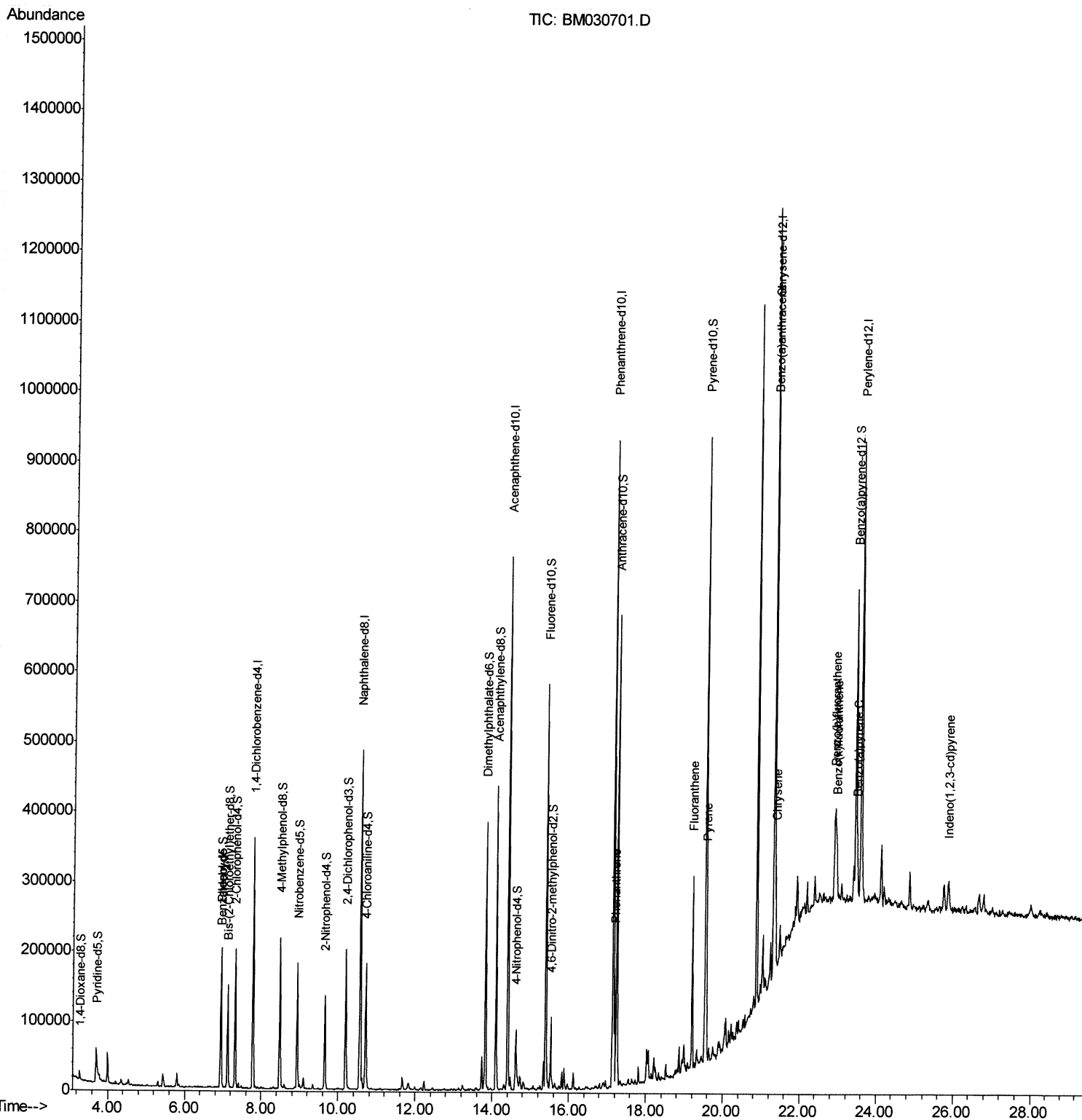
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM063021\
Data File : BM030701.D
Acq On : 30 Jun 2021 12:37
Operator : CG/JU
Sample : M2888-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB8H9

Quant Time: Jun 30 13:20:55 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 30 12:31:14 2021
Response via : Initial Calibration

Manual Integrations
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mohammad
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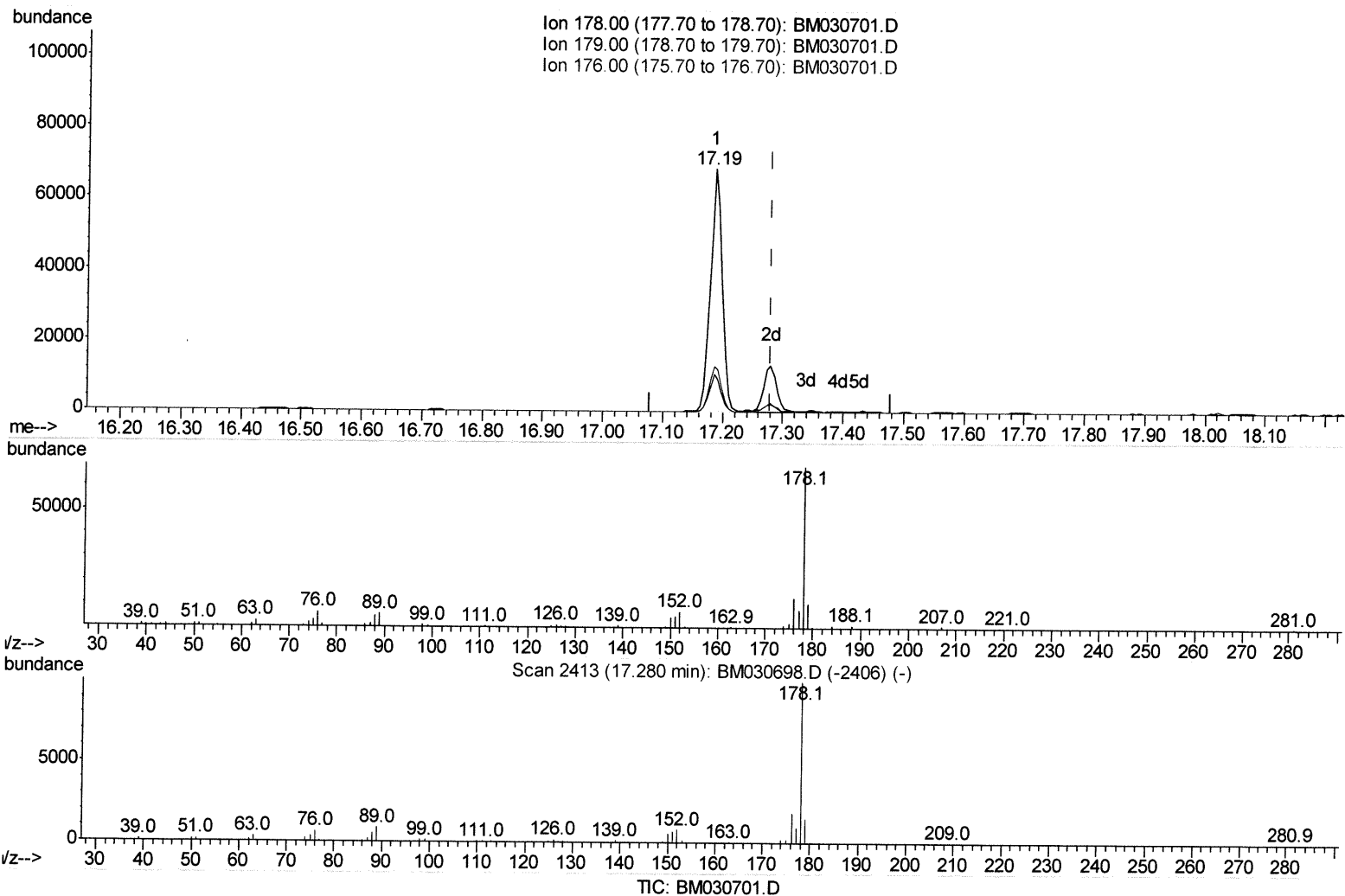
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(74) Anthracene

17.186min (-0.094) 3.16ng/ul

response 92958

Ion	Exp%	Act%
178.00	100	100
179.00	15.10	15.49
176.00	19.00	18.83
0.00	0.00	0.00

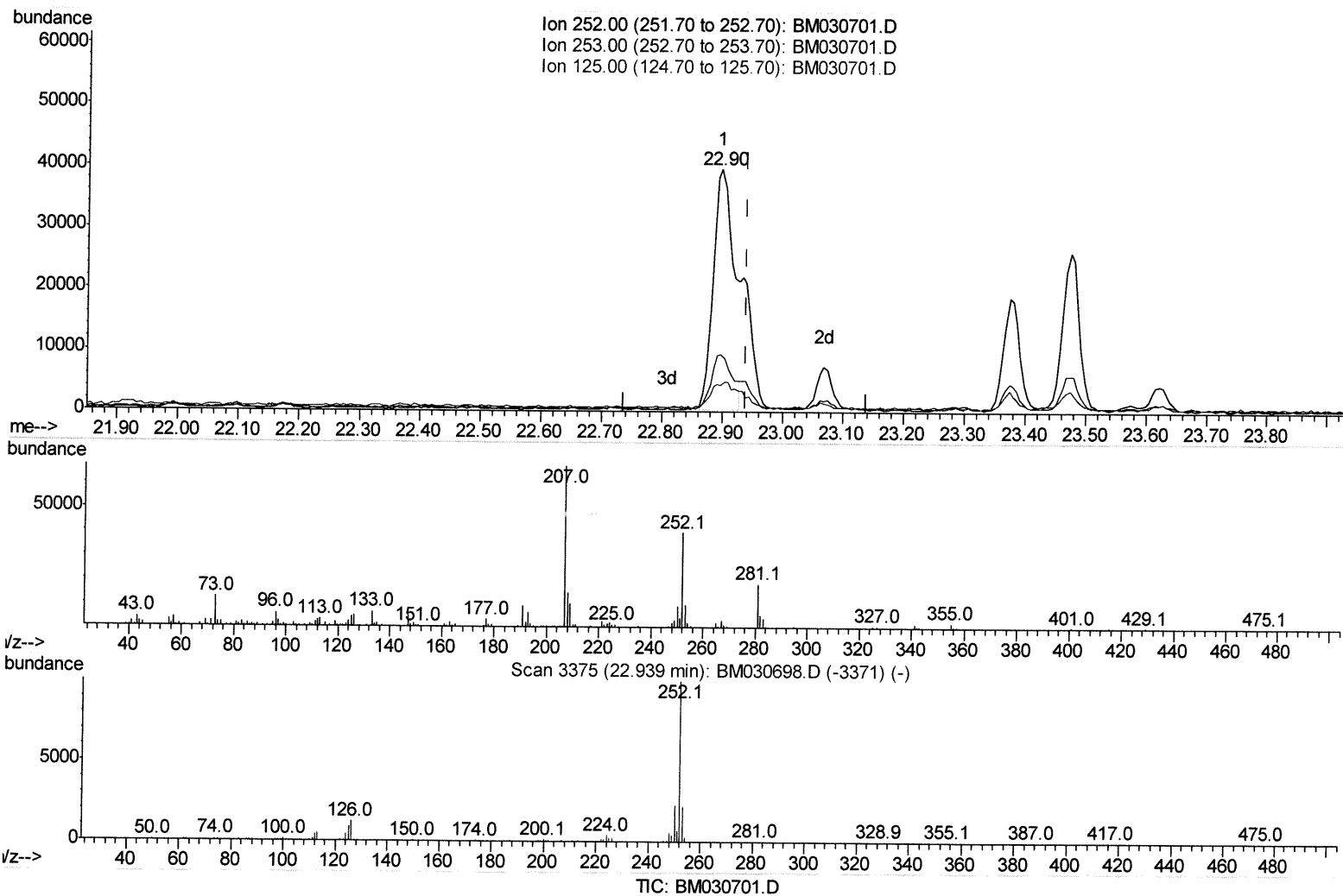
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(91) Benzo(k)fluoranthene

22.897min (-0.041) 3.40ng/ul

response 92101

Ion	Exp%	Act%
252.00	100	100
253.00	21.80	23.82
125.00	9.90	10.92
0.00	0.00	0.00

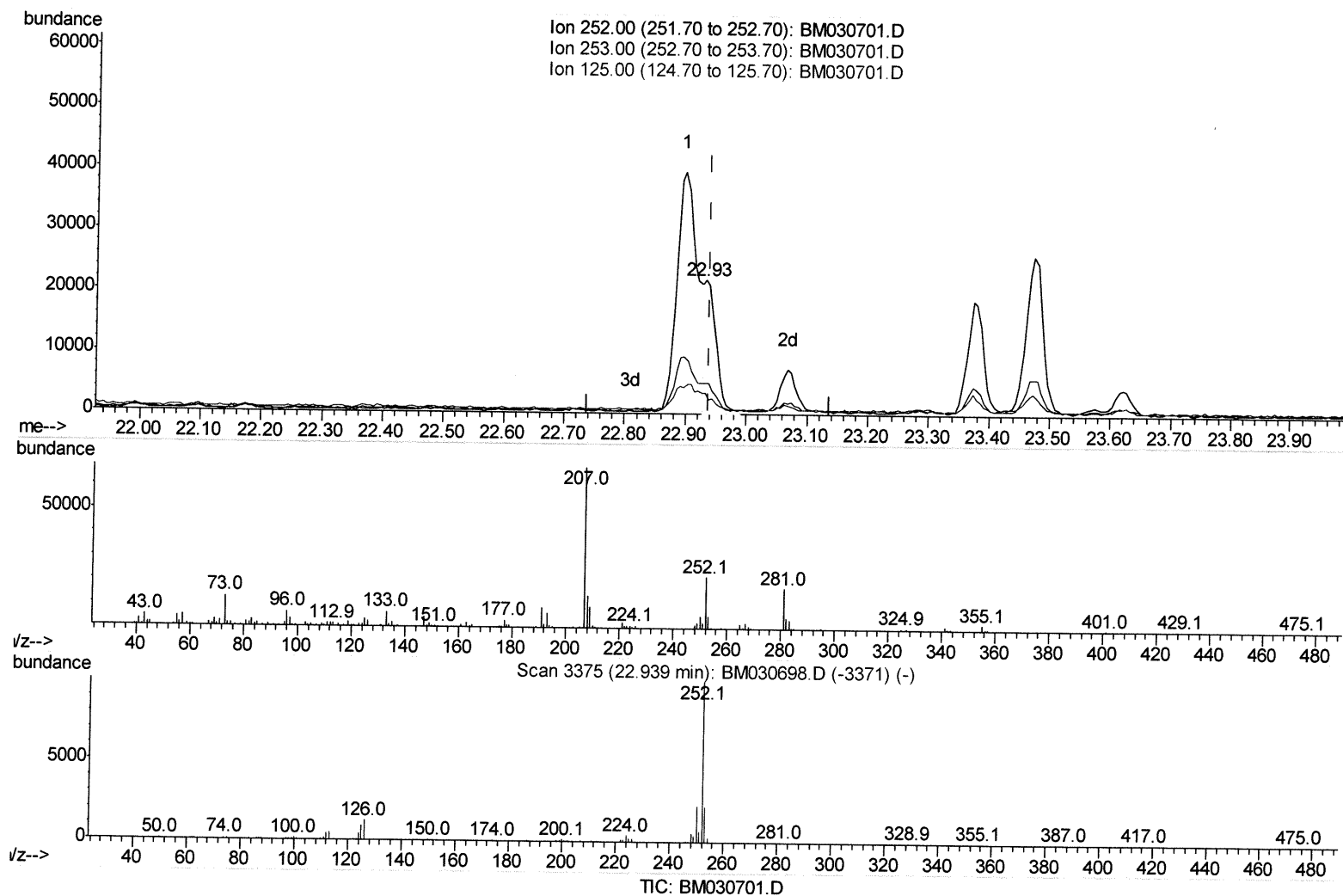
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(91) Benzo(k)fluoranthene

22.933min (-0.006) 1.09ng/ul m

response 29586

5/6/21

Ion	Exp%	Act%
252.00	100	100
253.00	21.80	23.17
125.00	9.90	15.79#
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	7.78	152	93675	20.00	ng/ul	0.00
20) Naphthalene-d8	10.56	136	388289	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.41	164	252788	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.14	188	536583	20.00	ng/ul	0.00
79) Chrysene-d12	21.32	240	504301	20.00	ng/ul	0.00
88) Perylene-d12	23.57	264	444302	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.27	96	6276	2.72	ng/uL	0.00
4) Pyridine-d5	3.70	84	39623	6.38	ng/ul	0.02
7) Phenol-d5	6.95	99	104281	12.91	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.12	67	69826	13.75	ng/ul	0.00
11) 2-Chlorophenol-d4	7.32	132	87328	14.00	ng/ul	0.00
15) 4-Methylphenol-d8	8.48	113	79037	12.57	ng/ul	0.00
21) Nitrobenzene-d5	8.93	128	39329	13.57	ng/ul	0.00
24) 2-Nitrophenol-d4	9.66	143	42267	13.13	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.19	165	78757	12.77	ng/ul	0.00
31) 4-Chloroaniline-d4	10.70	131	107266	12.47	ng/ul	0.00
46) Dimethylphthalate-d6	13.82	166	251437	14.41	ng/ul	0.00
49) Acenaphthylene-d8	14.10	160	308161	13.60	ng/ul	0.00
54) 4-Nitrophenol-d4	14.62	143	26594	8.04	ng/ul	0.00
60) Fluorene-d10	15.40	176	228383	14.75	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.53	200	23708	6.75	ng/ul	0.00
73) Anthracene-d10	17.24	188	357099	14.26	ng/ul	0.00
81) Pyrene-d10	19.53	212	405923	15.36	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.43	264	316814	13.72	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Benzaldehyde	6.93	77	4709	1.199	ng/ul	96
72) Phenanthrene	17.19	178	92958	3.226	ng/ul	99
80) Fluoranthene	19.20	202	159084	4.929	ng/ul	99
82) Pyrene	19.56	202	143398	4.229	ng/ul	99
85) Benzo(a)anthracene	21.30	228	82542	2.637	ng/ul	98
87) Chrysene	21.36	228	77241	2.532	ng/ul	99
90) Benzo(b)fluoranthene	22.90	252	92101	3.265	ng/ul	97
91) Benzo(k)fluoranthene	22.93	252	29586m>	1.091	ng/ul	> 96
93) Benzo(a)pyrene	23.47	252	49630	1.957	ng/ul#	96
94) Indeno(1,2,3-cd)pyrene	25.84	276	34895	1.093	ng/ul#	92

Jul 07/01/21

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