

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060220\
 Data File : BN010741.D
 Acq On : 02 Jun 2020 12:04
 Operator : CG/JU
 Sample : SST00.105
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.105

Quant Time: Jun 02 14:34:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 14:30:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1781	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5887	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3079	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6091	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	5314	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	5138	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	925	0.09	ng/ul	0.01
14) Fluoranthene-d10	18.98	212	1761	0.09	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	1802	0.096	ng/ul#	91
5) 2-Methylnaphthalene	12.00	142	1142	0.091	ng/ul	98
7) Acenaphthylene	13.91	152	1425	0.087	ng/ul#	92
8) Acenaphthene	14.26	153	1215	0.097	ng/ul	97
9) Fluorene	15.25	166	1281	0.086	ng/ul#	98
11) Pentachlorophenol	16.61	266	128	0.108	ng/ul#	86
12) Phenanthrene	16.98	178	2069	0.094	ng/ul	97
13) Anthracene	17.08	178	1623	0.084	ng/ul#	93
15) Fluoranthene	19.01	202	2218	0.083	ng/ul	93
17) Pyrene	19.37	202	2289	0.091	ng/ul#	92
18) Benzo(a)anthracene	21.13	228	1994	0.087	ng/ul	94
19) Chrysene	21.18	228	2519	0.096	ng/ul	98
21) Benzo(b)fluoranthene	22.66	252	2276	0.098	ng/ul	68
22) Benzo(k)fluoranthene	22.71	252	2477	0.101	ng/ul#	64
23) Benzo(a)pyrene	23.21	252	1859	0.092	ng/ul#	54
24) Indeno(1,2,3-cd)pyrene	25.42	276	2622	0.104	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.44	278	2077	0.100	ng/ul#	77
26) Benzo(g,h,i)perylene	26.06	276	2333	0.105	ng/ul#	90

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

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