

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098054.D
 Acq On : 25 Oct 2018 13:38
 Operator : SJ/JU
 Sample : SSTD1.665
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.665

Quant Time: Oct 25 14:13:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 25 13:24:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	923	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4782	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3343	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	7994	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10921	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9421	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	13660	1.82	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	42941	1.65	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	18462	1.604	ng/ul	97
5) 2-Methylnaphthalene	12.34	142	14070	1.839	ng/ul	99
7) Acenaphthylene	14.25	152	21627	1.836	ng/ul	97
8) Acenaphthene	14.60	153	15778	1.608	ng/ul	99
9) Fluorene	15.59	166	19142	1.618	ng/ul	99
11) Pentachlorophenol	16.95	266	715	0.980	ng/ul	96
12) Phenanthrene	17.33	178	30237	1.583	ng/ul	99
13) Anthracene	17.42	178	30644	1.889	ng/ul	98
15) Fluoranthene	19.35	202	38860	1.543	ng/ul	99
17) Pyrene	19.71	202	46913	1.748	ng/ul	98
18) Benzo(a)anthracene	21.45	228	46773	1.543	ng/ul	100
19) Chrysene	21.51	228	42454	1.452	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	43393	1.782	ng/ul	96
22) Benzo(k)fluoranthene	23.29	252	40824	1.479	ng/ul	96
23) Benzo(a)pyrene	23.91	252	41001	1.841	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.71	276	47537	1.880	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.73	278	39587	1.873	ng/ul#	96
26) Benzo(g,h,i)perylene	27.54	276	39749	1.665	ng/ul	96

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

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