

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096957.D
 Acq On : 17 Jul 2018 13:37
 Operator : SJ/JU
 Sample : SSTD1.665
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.665

Quant Time: Jul 17 14:16:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 13:01:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1503	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6948	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3687	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	9949	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	13708	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11795	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	18161	1.60	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	52607	2.12	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.19	128	27452	1.635	ng/ul#	90
5) 2-Methylnaphthalene	11.81	142	18528	1.620	ng/ul	97
7) Acenaphthylene	13.74	152	26618	1.761	ng/ul	95
8) Acenaphthene	14.08	153	20447	1.639	ng/ul	93
9) Fluorene	15.08	166	22327	1.734	ng/ul	92
11) Pentachlorophenol	16.43	266	1936	2.973	ng/ul	97
12) Phenanthrene	16.81	178	42762	1.667	ng/ul#	88
13) Anthracene	16.90	178	39816	1.828	ng/ul#	91
15) Fluoranthene	18.84	202	58799	2.152	ng/ul	84
17) Pyrene	19.21	202	57365	1.275	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	60464	1.870	ng/ul#	86
19) Chrysene	21.01	228	63713	1.521	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	58771	1.826	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	59717	1.699	ng/ul#	83
23) Benzo(a)pyrene	23.11	252	52143	1.915	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	59341	2.099	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	50026	2.228	ng/ul#	86
26) Benzo(g,h,i)perylene	26.20	276	51233	1.867	ng/ul#	92

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

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