

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097854.D
 Acq On : 11 Oct 2018 12:51
 Operator : SJ/JU
 Sample : SSTD1.648
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.648

Quant Time: Oct 11 13:29:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 12:14:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2284	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10792	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7723	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	23459	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	29076	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	27589	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	30075	1.78	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	128982	1.69	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	42580	1.639	ng/ul	100
5) 2-Methylnaphthalene	12.36	142	31522	1.825	ng/ul	99
7) Acenaphthylene	14.26	152	49508	1.819	ng/ul	98
8) Acenaphthene	14.60	153	37481	1.653	ng/ul	97
9) Fluorene	15.59	166	49701	1.819	ng/ul	99
11) Pentachlorophenol	16.94	266	5829	2.723	ng/ul	98
12) Phenanthrene	17.34	178	91933	1.640	ng/ul	98
13) Anthracene	17.43	178	88614	1.862	ng/ul	98
15) Fluoranthene	19.36	202	123030	1.665	ng/ul	98
17) Pyrene	19.72	202	149808	2.097	ng/ul	98
18) Benzo(a)anthracene	21.46	228	143170	1.774	ng/ul	98
19) Chrysene	21.51	228	135713	1.743	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	126986	1.781	ng/ul	98
22) Benzo(k)fluoranthene	23.30	252	128761	1.593	ng/ul#	96
23) Benzo(a)pyrene	23.91	252	121122	1.857	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.71	276	139077	1.878	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	115802	1.871	ng/ul	99
26) Benzo(g,h,i)perylene	27.54	276	118594	1.696	ng/ul	99

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

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