

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042518\
 Data File : BE096164.D
 Acq On : 26 Apr 2018 00:54
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.463

Quant Time: Apr 26 07:16:46 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 26 06:40:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	734	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2859	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1856	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	5031	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	6508	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	5079	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2194	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	7861	0.34	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	11.23	128	3180	0.399	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	2258	0.423	ng/ul	99
7) Acenaphthylene	14.67	152	3796	0.397	ng/ul	98
8) Acenaphthene	15.01	153	2832	0.411	ng/ul	96
9) Fluorene	15.97	166	4047	0.443	ng/ul#	89
11) Pentachlorophenol	17.32	266	332	0.274	ng/ul	95
12) Phenanthrene	17.70	178	6279	0.424	ng/ul#	89
13) Anthracene	17.79	178	5923	0.399	ng/ul#	91
15) Fluoranthene	19.67	202	7865	0.347	ng/ul	85
17) Pyrene	20.02	202	4368	0.335	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	5841	0.296	ng/ul#	86
19) Chrysene	21.79	228	6820	0.377	ng/ul	96
21) Benzo(b)fluoranthene	23.55	252	6957	0.362	ng/ul	86
22) Benzo(k)fluoranthene	23.59	252	6441	0.350	ng/ul#	90
23) Benzo(a)pyrene	24.23	252	6391	0.383	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.13	276	7369	0.327	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.14	278	6053	0.321	ng/ul	92
26) Benzo(g,h,i)perylene	27.99	276	6098	0.311	ng/ul	96

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

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