

Data Path : Z:\HPCHEM1\BNA_E\Data\BE042518\
 Data File : BE096141.D
 Acq On : 25 Apr 2018 9:07
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.461

Quant Time: Apr 26 04:14:41 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 25 05:27:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	740	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2880	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1587	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4495	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	6744	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	5268	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2083	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	7958	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.23	128	3165	0.395	ng/ul#	91
5) 2-Methylnaphthalene	12.81	142	2114	0.393	ng/ul	99
7) Acenaphthylene	14.68	152	3269	0.400	ng/ul	98
8) Acenaphthene	15.01	153	2423	0.411	ng/ul	93
9) Fluorene	15.98	166	2753	0.353	ng/ul	94
11) Pentachlorophenol	17.31	266	352	0.325	ng/ul	97
12) Phenanthrene	17.70	178	5280	0.399	ng/ul#	89
13) Anthracene	17.79	178	5194	0.391	ng/ul#	92
15) Fluoranthene	19.67	202	7874	0.389	ng/ul	84
17) Pyrene	20.02	202	4384	0.324	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	5832	0.286	ng/ul#	86
19) Chrysene	21.79	228	7225	0.385	ng/ul	96
21) Benzo(b)fluoranthene	23.54	252	7175	0.360	ng/ul	87
22) Benzo(k)fluoranthene	23.60	252	6778	0.355	ng/ul#	88
23) Benzo(a)pyrene	24.23	252	6602	0.381	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.13	276	7567	0.324	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.15	278	6175	0.316	ng/ul	90
26) Benzo(g,h,i)perylene	28.00	276	6287	0.309	ng/ul	95

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

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