

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042518\
 Data File : BE096171.D
 Acq On : 26 Apr 2018 6:28
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.464

Quant Time: Apr 26 07:14:07 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	8.35	152	622	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2719	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1612	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3623	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	4693	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3706	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2073	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5594	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.23	128	3014	0.398	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	2107	0.415	ng/ul	96
7) Acenaphthylene	14.67	152	3352	0.403	ng/ul	98
8) Acenaphthene	15.01	153	2521	0.421	ng/ul	92
9) Fluorene	15.98	166	3225	0.407	ng/ul	93
11) Pentachlorophenol	17.32	266	232	0.266	ng/ul	92
12) Phenanthrene	17.70	178	5175	0.485	ng/ul#	89
13) Anthracene	17.79	178	4439	0.415	ng/ul#	92
15) Fluoranthene	19.67	202	6232	0.382	ng/ul	85
17) Pyrene	20.02	202	4130	0.439	ng/ul#	80
18) Benzo(a)anthracene	21.73	228	4438	0.312	ng/ul#	86
19) Chrysene	21.79	228	5116	0.392	ng/ul	96
21) Benzo(b)fluoranthene	23.55	252	5156	0.367	ng/ul	91
22) Benzo(k)fluoranthene	23.60	252	4879	0.364	ng/ul	92
23) Benzo(a)pyrene	24.24	252	4705	0.386	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	27.14	276	5323	0.324	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.15	278	4327	0.315	ng/ul	95
26) Benzo(g,h,i)perylene	28.01	276	4487	0.314	ng/ul	97

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

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