

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE071118\
 Data File : BE096895.D
 Acq On : 12 Jul 2018 12:09
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.457

Quant Time: Jul 12 14:23:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 10 02:37:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2682	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	14936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	8204	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	19354	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	16685	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	16837	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8360	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15611	0.32	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	12170	0.337	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	8475	0.345	ng/ul	98
7) Acenaphthylene	13.74	152	11629	0.346	ng/ul	96
8) Acenaphthene	14.09	153	8996	0.324	ng/ul#	95
9) Fluorene	15.08	166	9010	0.315	ng/ul#	93
11) Pentachlorophenol	16.43	266	431	0.340	ng/ul	98
12) Phenanthrene	16.82	178	16289	0.326	ng/ul#	89
13) Anthracene	16.91	178	14015	0.331	ng/ul#	92
15) Fluoranthene	18.85	202	17365	0.327	ng/ul	84
17) Pyrene	19.21	202	16812	0.307	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	13704	0.348	ng/ul#	85
19) Chrysene	21.02	228	16094	0.316	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	14245	0.310	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	14547	0.290	ng/ul#	86
23) Benzo(a)pyrene	23.12	252	12399	0.319	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.51	276	13987	0.347	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.54	278	11440	0.357	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	12312	0.314	ng/ul#	93

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

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