

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE070918\
 Data File : BE096889.D
 Acq On : 9 Jul 2018 20:02
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.454

Quant Time: Jul 10 03:14:53 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	2774	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	15423	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	9136	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	21759	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	21324	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	22769	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8795	0.35	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	19842	0.37	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.19	128	12501	0.335	ng/ul#	91
5) 2-Methylnaphthalene	11.82	142	8837	0.348	ng/ul	97
7) Acenaphthylene	13.74	152	12987	0.347	ng/ul	96
8) Acenaphthene	14.09	153	9787	0.317	ng/ul	94
9) Fluorene	15.08	166	10276	0.322	ng/ul	91
11) Pentachlorophenol	16.43	266	954	0.670	ng/ul	96
12) Phenanthrene	16.82	178	17786	0.317	ng/ul#	89
13) Anthracene	16.91	178	17002	0.357	ng/ul#	92
15) Fluoranthene	18.85	202	20895	0.350	ng/ul	84
17) Pyrene	19.21	202	21795	0.311	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	20262	0.403	ng/ul#	84
19) Chrysene	21.02	228	19946	0.306	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	21774	0.350	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	20505	0.302	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	20139	0.383	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	22620	0.415	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	18435	0.425	ng/ul#	89
26) Benzo(g,h,i)perylene	26.20	276	18918	0.357	ng/ul#	93

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

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