

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096857.D
 Acq On : 6 Jul 2018 17:29
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.451

Quant Time: Jul 06 18:07:15 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 : Fri Jul 06 16:55:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2950	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	14627	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	8007	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	21599	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	17775	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	14638	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	9574	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	21035	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.19	128	14294	0.404	ng/ul#	90
5) 2-Methylnaphthalene	11.81	142	9665	0.401	ng/ul	99
7) Acenaphthylene	13.74	152	13130	0.400	ng/ul	96
8) Acenaphthene	14.09	153	10793	0.398	ng/ul#	95
9) Fluorene	15.08	166	11186	0.400	ng/ul#	92
11) Pentachlorophenol	16.43	266	494	0.349	ng/ul	97
12) Phenanthrene	16.82	178	22285	0.400	ng/ul#	89
13) Anthracene	16.90	178	18508	0.391	ng/ul#	92
15) Fluoranthene	18.85	202	23518	0.396	ng/ul	84
17) Pyrene	19.21	202	23985	0.411	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	16628	0.397	ng/ul#	84
19) Chrysene	21.02	228	21715	0.400	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	15963	0.400	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	17635	0.404	ng/ul#	86
23) Benzo(a)pyrene	23.12	252	13483	0.399	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.51	276	14108	0.402	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.53	278	11334	0.407	ng/ul#	90
26) Benzo(g,h,i)perylene	26.21	276	13738	0.403	ng/ul#	93

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

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