

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096853.D
 Acq On : 6 Jul 2018 14:12
 Operator : SJ/JU
 Sample : SST0.847
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.847

Quant Time: Jul 06 14:58:54 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 14:58:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2915	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	14945	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	8169	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	21776	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	18547	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	14776	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	19351	0.75	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	43574	0.68	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	28378	0.744	ng/ul#	90
5) 2-Methylnaphthalene	11.81	142	19821	0.801	ng/ul	97
7) Acenaphthylene	13.74	152	26890	0.758	ng/ul	96
8) Acenaphthene	14.09	153	22024	0.751	ng/ul	94
9) Fluorene	15.08	166	23168	0.738	ng/ul#	92
11) Pentachlorophenol	16.43	266	1063	1.255	ng/ul	97
12) Phenanthrene	16.82	178	44341	0.726	ng/ul#	89
13) Anthracene	16.91	178	38691	0.762	ng/ul#	92
15) Fluoranthene	18.85	202	48749	0.721	ng/ul	84
17) Pyrene	19.21	202	45518	0.701	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	35413	0.796	ng/ul#	85
19) Chrysene	21.02	228	43963	0.707	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	34067	0.781	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	37679	0.830	ng/ul#	84
23) Benzo(a)pyrene	23.12	252	29011	0.741	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.51	276	30607	0.778	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.54	278	24964	0.821	ng/ul#	87
26) Benzo(g,h,i)perylene	26.20	276	29112	0.746	ng/ul#	92

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

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