

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096955.D
 Acq On : 17 Jul 2018 12:19
 Operator : SJ/JU
 Sample : SST0.463
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.463

Quant Time: Jul 17 13:01:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 13:01:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1678	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	7736	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3886	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	10500	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12970	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12149	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4856	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12328	0.47	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	7463	0.399	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	4856	0.381	ng/ul	98
7) Acenaphthylene	13.74	152	6747	0.424	ng/ul	96
8) Acenaphthene	14.09	153	5199	0.395	ng/ul	95
9) Fluorene	15.08	166	5410	0.399	ng/ul	92
11) Pentachlorophenol	16.43	266	454	0.661	ng/ul	98
12) Phenanthrene	16.81	178	10929	0.404	ng/ul#	88
13) Anthracene	16.90	178	9448	0.411	ng/ul#	91
15) Fluoranthene	18.84	202	13915	0.483	ng/ul	84
17) Pyrene	19.21	202	14583	0.343	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	13387	0.438	ng/ul#	86
19) Chrysene	21.01	228	15168	0.383	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	13408	0.404	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	13922	0.385	ng/ul#	84
23) Benzo(a)pyrene	23.11	252	11669	0.416	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	13300	0.457	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	11178	0.483	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	11863	0.420	ng/ul#	93

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

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