

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096956.D
 Acq On : 17 Jul 2018 12:58
 Operator : SJ/JU
 Sample : SSTD0.864
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.864

Quant Time: Jul 17 13:48:12 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	7.39	152	1487	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6868	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3567	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10257	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	13003	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11370	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8764	0.78	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	24908	0.98	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	13160	0.793	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	8768	0.775	ng/ul	97
7) Acenaphthylene	13.74	152	12439	0.851	ng/ul	96
8) Acenaphthene	14.08	153	9576	0.793	ng/ul#	95
9) Fluorene	15.08	166	10483	0.842	ng/ul	94
11) Pentachlorophenol	16.43	266	858	1.278	ng/ul	96
12) Phenanthrene	16.81	178	20940	0.792	ng/ul#	88
13) Anthracene	16.90	178	18781	0.836	ng/ul#	91
15) Fluoranthene	18.84	202	28064	0.996	ng/ul	84
17) Pyrene	19.21	202	27602	0.647	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	27197	0.887	ng/ul#	86
19) Chrysene	21.01	228	29793	0.750	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	26931	0.868	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	26388	0.779	ng/ul#	84
23) Benzo(a)pyrene	23.11	252	23014	0.877	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.51	276	26279	0.964	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	21973	1.015	ng/ul#	87
26) Benzo(g,h,i)perylene	26.20	276	23086	0.873	ng/ul#	93

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

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