

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096952.D
 Acq On : 17 Jul 2018 10:26
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.401

Quant Time: Jul 18 10:45:50 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 : Wed Jul 18 07:15:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1590	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	7309	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3720	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	10518	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	13878	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	13590	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4580	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12872	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.19	128	7045	0.398	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	4658	0.398	ng/ul	98
7) Acenaphthylene	13.74	152	6574	0.405	ng/ul	97
8) Acenaphthene	14.09	153	5018	0.399	ng/ul	94
9) Fluorene	15.08	166	5341	0.399	ng/ul	94
11) Pentachlorophenol	16.43	266	491	0.413	ng/ul	96
12) Phenanthrene	16.81	178	10798	0.391	ng/ul#	87
13) Anthracene	16.90	178	9535	0.395	ng/ul#	91
15) Fluoranthene	18.84	202	14478	0.398	ng/ul	84
17) Pyrene	19.21	202	15647	0.390	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	14870	0.405	ng/ul#	86
19) Chrysene	21.01	228	16097	0.386	ng/ul	95
21) Benzo(b)fluoranthene	22.54	252	15321	0.397	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	15235	0.382	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	13570	0.401	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	15370	0.405	ng/ul#	80
25) Dibenzo(a,h)anthracene	25.53	278	12679	0.400	ng/ul#	88
26) Benzo(g,h,i)perylene	26.20	276	13518	0.399	ng/ul	93

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

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