

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE071818\
 Data File : BE096998.D
 Acq On : 18 Jul 2018 17:55
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.472

Quant Time: Jul 18 18:33:06 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 15:28:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1384	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6840	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	4359	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	12307	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	13683	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	13319	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4591	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15224	0.41	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.19	128	6621	0.399	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	4654	0.425	ng/ul	97
7) Acenaphthylene	13.74	152	7449	0.392	ng/ul	96
8) Acenaphthene	14.09	153	5510	0.374	ng/ul	94
9) Fluorene	15.08	166	6580	0.420	ng/ul	93
11) Pentachlorophenol	16.43	266	651	0.468	ng/ul	96
12) Phenanthrene	16.81	178	12711	0.393	ng/ul#	88
13) Anthracene	16.90	178	11726	0.416	ng/ul#	90
15) Fluoranthene	18.84	202	15338	0.360	ng/ul	84
17) Pyrene	19.21	202	15794	0.399	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	15132	0.418	ng/ul#	85
19) Chrysene	21.02	228	15666	0.381	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	14957	0.396	ng/ul	84
22) Benzo(k)fluoranthene	22.58	252	14177	0.363	ng/ul#	86
23) Benzo(a)pyrene	23.11	252	13349	0.402	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	16782	0.451	ng/ul#	77
25) Dibenzo(a,h)anthracene	25.53	278	13942	0.449	ng/ul#	86
26) Benzo(g,h,i)perylene	26.21	276	14558	0.439	ng/ul#	91

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

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