

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098073.D
 Acq On : 26 Oct 2018 2:00
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.469

Quant Time: Oct 26 05:38:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 26 05:34:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1083	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5379	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	3703	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8847	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	11851	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	10122	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3562	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10647	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.73	128	5032	0.391	ng/ul	99
5) 2-Methylnaphthalene	12.35	142	3682	0.378	ng/ul	99
7) Acenaphthylene	14.25	152	5557	0.380	ng/ul	98
8) Acenaphthene	14.60	153	4066	0.372	ng/ul	99
9) Fluorene	15.59	166	4863	0.370	ng/ul	98
11) Pentachlorophenol	16.96	266	239	0.516	ng/ul	96
12) Phenanthrene	17.33	178	7717	0.378	ng/ul	98
13) Anthracene	17.42	178	7703	0.382	ng/ul	99
15) Fluoranthene	19.35	202	9633	0.370	ng/ul	99
17) Pyrene	19.72	202	9765	0.353	ng/ul	99
18) Benzo(a)anthracene	21.46	228	13880	0.398	ng/ul	98
19) Chrysene	21.51	228	10342	0.384	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	10383	0.370	ng/ul	97
22) Benzo(k)fluoranthene	23.29	252	10560	0.385	ng/ul	97
23) Benzo(a)pyrene	23.91	252	10088	0.374	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.72	276	11754	0.380	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.75	278	9877	0.383	ng/ul#	95
26) Benzo(g,h,i)perylene	27.55	276	9852	0.379	ng/ul	97

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

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