

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098106.D
 Acq On : 27 Oct 2018 1:52
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.473

Quant Time: Oct 27 04:04:59 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 : Sat Oct 27 00:33:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1059	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5336	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3588	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8806	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	12840	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	10982	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3576	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10449	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.72	128	4866	0.381	ng/ul	100
5) 2-Methylnaphthalene	12.34	142	3626	0.376	ng/ul	100
7) Acenaphthylene	14.25	152	5401	0.382	ng/ul	99
8) Acenaphthene	14.60	153	4023	0.380	ng/ul	99
9) Fluorene	15.59	166	4739	0.372	ng/ul	96
11) Pentachlorophenol	16.95	266	359	0.779	ng/ul	98
12) Phenanthrene	17.33	178	7535	0.371	ng/ul	99
13) Anthracene	17.42	178	7646	0.381	ng/ul	100
15) Fluoranthene	19.35	202	9530	0.368	ng/ul	98
17) Pyrene	19.71	202	9617	0.321	ng/ul	99
18) Benzo(a)anthracene	21.45	228	15141	0.401	ng/ul	99
19) Chrysene	21.51	228	10364	0.355	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	10653	0.350	ng/ul	98
22) Benzo(k)fluoranthene	23.29	252	9978	0.335	ng/ul#	99
23) Benzo(a)pyrene	23.91	252	10082	0.345	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.72	276	11854	0.353	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.74	278	9751	0.349	ng/ul	97
26) Benzo(g,h,i)perylene	27.55	276	9805	0.347	ng/ul	97

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

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