

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097857.D
 Acq On : 11 Oct 2018 14:59
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.451

Quant Time: Oct 11 16:30:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 14:16:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2347	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11009	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7841	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	23902	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	28478	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	28642	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7361	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	30165	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.74	128	10602	0.396	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	7514	0.388	ng/ul	98
7) Acenaphthylene	14.26	152	11829	0.399	ng/ul	99
8) Acenaphthene	14.60	153	9257	0.402	ng/ul	98
9) Fluorene	15.59	166	11969	0.398	ng/ul	98
11) Pentachlorophenol	16.94	266	1154	0.342	ng/ul	99
12) Phenanthrene	17.34	178	22647	0.400	ng/ul	99
13) Anthracene	17.43	178	20408	0.389	ng/ul	99
15) Fluoranthene	19.36	202	28745	0.396	ng/ul	99
17) Pyrene	19.72	202	35266	0.416	ng/ul	99
18) Benzo(a)anthracene	21.47	228	36645	0.426	ng/ul	99
19) Chrysene	21.51	228	32567	0.421	ng/ul	97
21) Benzo(b)fluoranthene	23.25	252	29608	0.387	ng/ul	99
22) Benzo(k)fluoranthene	23.30	252	32267	0.394	ng/ul	97
23) Benzo(a)pyrene	23.91	252	28739	0.382	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.72	276	32875	0.385	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	27482	0.390	ng/ul	98
26) Benzo(g,h,i)perylene	27.54	276	28969	0.390	ng/ul	98

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

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