

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080118\
 Data File : BE097130.D
 Acq On : 1 Aug 2018 14:56
 Operator : SJ/JU
 Sample : SSTD1.677
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.677

Quant Time: Aug 01 15:38:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 01 14:22:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1427	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	6830	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3897	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10127	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12884	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11880	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	18205	1.63	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	56582	2.24	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	26649	1.614	ng/ul#	91
5) 2-Methylnaphthalene	11.80	142	18088	1.609	ng/ul	98
7) Acenaphthylene	13.73	152	27322	1.711	ng/ul	95
8) Acenaphthene	14.08	153	19826	1.503	ng/ul	93
9) Fluorene	15.08	166	23098	1.698	ng/ul#	92
11) Pentachlorophenol	16.43	266	2549	3.846	ng/ul	98
12) Phenanthrene	16.81	178	42193	1.616	ng/ul#	89
13) Anthracene	16.90	178	40837	1.842	ng/ul#	92
15) Fluoranthene	18.84	202	55871	2.009	ng/ul	84
17) Pyrene	19.20	202	55804	1.319	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	59588	1.961	ng/ul#	84
19) Chrysene	21.01	228	60963	1.548	ng/ul	97
21) Benzo(b)fluoranthene	22.53	252	55325	1.706	ng/ul	81
22) Benzo(k)fluoranthene	22.58	252	58006	1.639	ng/ul#	84
23) Benzo(a)pyrene	23.11	252	53047	1.934	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	68372	2.402	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.52	278	56359	2.492	ng/ul#	85
26) Benzo(g,h,i)perylene	26.19	276	59098	2.139	ng/ul#	90

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

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