

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094767.D
 Acq On : 6 Nov 2017 23:22
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.479

Manual Integrations
 APPROVED

Sohil
 11/7/2017 5:42:35 PM

Quant Time: Nov 07 01:53:46 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 01:23:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1101	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6556	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3631	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	7937	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7853	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	7365	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	3913	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9000	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	5687	0.355	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	3936	0.351	ng/ul	99
7) Acenaphthylene	14.25	152	6065	0.369	ng/ul	98
8) Acenaphthene	14.58	153	4262	0.361	ng/ul	95
9) Fluorene	15.57	166	4838	0.365	ng/ul	91
11) Pentachlorophenol	17.00	266	175m	0.442	ng/ul	
12) Phenanthrene	17.31	178	7580	0.362	ng/ul	93
13) Anthracene	17.40	178	8286m	0.401	ng/ul	
15) Fluoranthene	19.32	202	9097	0.375	ng/ul	84
17) Pyrene	19.68	202	9758	0.402	ng/ul#	81
18) Benzo(a)anthracene	21.42	228	8944	0.398	ng/ul#	87
19) Chrysene	21.47	228	8317	0.358	ng/ul#	84
21) Benzo(b)fluoranthene	23.20	252	8109	0.391	ng/ul#	81
22) Benzo(k)fluoranthene	23.25	252	8702	0.392	ng/ul#	77
23) Benzo(a)pyrene	23.87	252	8161	0.394	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.70	276	7972	0.372	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.69	278	6736	0.373	ng/ul#	66
26) Benzo(g,h,i)perylene	27.55	276	6571	0.382	ng/ul#	69

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

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