

Data Path : Z:\HPCHEM1\BNA_E\Data\BE082917\
 Data File : BE093899.D
 Acq On : 28 Aug 2017 16:59
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.415

Quant Time: Aug 28 17:40:44 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 28 15:03:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2028	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	10591	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5884	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	13263	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13995	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	13156	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	6698	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	16192	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	10400	0.404	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	7086	0.377	ng/ul	99
7) Acenaphthylene	14.24	152	10684	0.434	ng/ul	96
8) Acenaphthene	14.59	153	7712	0.398	ng/ul	98
9) Fluorene	15.57	166	8859	0.362	ng/ul#	90
11) Pentachlorophenol	16.93	266	951	0.435	ng/ul	96
12) Phenanthrene	17.29	178	14275	0.395	ng/ul#	89
13) Anthracene	17.38	178	14483	0.412	ng/ul#	89
15) Fluoranthene	19.30	202	16781	0.372	ng/ul	84
17) Pyrene	19.67	202	18106	0.443	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	16147	0.398	ng/ul#	87
19) Chrysene	21.45	228	16050	0.412	ng/ul	96
21) Benzo(b)fluoranthene	23.14	252	15402	0.378	ng/ul	91
22) Benzo(k)fluoranthene	23.19	252	15387	0.383	ng/ul	93
23) Benzo(a)pyrene	23.80	252	15019	0.388	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.56	276	15692	0.367	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.59	278	12761	0.357	ng/ul	93
26) Benzo(g,h,i)perylene	27.39	276	13854	0.385	ng/ul	94

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

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