

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061517\
 Data File : BE093241.D
 Acq On : 16 Jun 2017 16:16
 Operator : SJ/MA
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.412

Quant Time: Jun 16 17:01:19 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 16:46:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2534	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13061	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	7393	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	17796	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	18368	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	15872	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	8326	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	20410	0.42	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.57	128	13453	0.415	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	9576	0.432	ng/ul	98
7) Acenaphthylene	14.09	152	14510	0.457	ng/ul	97
8) Acenaphthene	14.43	153	10597	0.414	ng/ul	98
9) Fluorene	15.42	166	11921	0.397	ng/ul	90
11) Pentachlorophenol	16.77	266	1325	0.502	ng/ul	98
12) Phenanthrene	17.13	178	19914	0.407	ng/ul#	87
13) Anthracene	17.22	178	19971	0.468	ng/ul#	86
15) Fluoranthene	19.16	202	24551	0.409	ng/ul	83
17) Pyrene	19.52	202	26108	0.482	ng/ul#	83
18) Benzo(a)anthracene	21.25	228	22648	0.462	ng/ul#	86
19) Chrysene	21.30	228	20976	0.413	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	20063	0.417	ng/ul	94
22) Benzo(k)fluoranthene	22.98	252	20011	0.391	ng/ul	95
23) Benzo(a)pyrene	23.55	252	19220	0.422	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.19	276	21901	0.415	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	18096	0.407	ng/ul#	89
26) Benzo(g,h,i)perylene	26.96	276	18087	0.400	ng/ul	94

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

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