

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062117\
 Data File : BE093308.D
 Acq On : 21 Jun 2017 9:44
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.419

Quant Time: Jun 22 02:08:33 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1819	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9751	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5529	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13340	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13497	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11404	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	6208	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	14368	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.56	128	9989	0.41	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	7181	0.43	ng/ul	99
7) Acenaphthylene	14.08	152	10288	0.43	ng/ul	97
8) Acenaphthene	14.42	153	7914	0.41	ng/ul	97
9) Fluorene	15.41	166	8924	0.40	ng/ul	90
11) Pentachlorophenol	16.75	266	790	0.40	ng/ul	96
12) Phenanthrene	17.13	178	15013	0.41	ng/ul#	89
13) Anthracene	17.22	178	14128	0.44	ng/ul#	89
15) Fluoranthene	19.15	202	17616	0.39	ng/ul	83
17) Pyrene	19.51	202	18942	0.48	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	16290	0.45	ng/ul#	86
19) Chrysene	21.29	228	15590	0.42	ng/ul	95
21) Benzo(b)fluoranthene	22.93	252	14679	0.42	ng/ul	88
22) Benzo(k)fluoranthene	22.97	252	14592	0.40	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	13944	0.43	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	15884	0.42	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.20	278	13205	0.41	ng/ul#	88
26) Benzo(g,h,i)perylene	26.96	276	13048	0.40	ng/ul	97

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

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