

Data Path : Z:\HPCHEM1\BNA_E\Data\BE060817\
 Data File : BE093073.D
 Acq On : 8 Jun 2017 8:58
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.458

Quant Time: Jun 08 13:24:54 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 13:24:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	4256	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	19167	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9580	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	25846	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	34643	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	30078	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	10821	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	29113	0.41	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	18841	0.396	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	12530	0.385	ng/ul	99
7) Acenaphthylene	14.09	152	15990	0.389	ng/ul	97
8) Acenaphthene	14.43	153	13345	0.402	ng/ul	96
9) Fluorene	15.42	166	15218	0.391	ng/ul#	92
11) Pentachlorophenol	16.75	266	1292	0.337	ng/ul	99
12) Phenanthrene	17.13	178	28004	0.394	ng/ul#	85
13) Anthracene	17.24	178	24838	0.401	ng/ul	92
15) Fluoranthene	19.16	202	35988	0.413	ng/ul	84
17) Pyrene	19.52	202	41173	0.403	ng/ul#	84
18) Benzo(a)anthracene	21.25	228	37756	0.409	ng/ul#	86
19) Chrysene	21.30	228	37972	0.396	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	33807	0.370	ng/ul	87
22) Benzo(k)fluoranthene	22.98	252	38644	0.399	ng/ul#	90
23) Benzo(a)pyrene	23.56	252	33921	0.393	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.19	276	39038	0.391	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.21	278	33200	0.395	ng/ul#	88
26) Benzo(g,h,i)perylene	26.96	276	33263	0.388	ng/ul	96

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

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