

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031318\
 Data File : BE095752.D
 Acq On : 13 Mar 2018 16:24
 Operator : SJ/JU
 Sample : SSTD1.613
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.613

Quant Time: Mar 13 17:04:09 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 13 16:24:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1774	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5004	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	2822	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6192	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6569	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	5753	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	14107	1.73	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	34098	1.44	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	22893	1.634	ng/ul#	87
5) 2-Methylnaphthalene	12.87	142	16207	1.732	ng/ul	100
7) Acenaphthylene	14.73	152	24397	1.796	ng/ul	96
8) Acenaphthene	15.07	153	17858	1.695	ng/ul	94
9) Fluorene	16.03	166	19744	1.818	ng/ul#	89
11) Pentachlorophenol	17.36	266	1645	1.372	ng/ul	98
12) Phenanthrene	17.75	178	31181	1.497	ng/ul#	88
13) Anthracene	17.84	178	30647	1.608	ng/ul#	91
15) Fluoranthene	19.72	202	39171	1.453	ng/ul	83
17) Pyrene	20.07	202	41000	1.860	ng/ul#	77
18) Benzo(a)anthracene	21.78	228	35569	1.732	ng/ul#	86
19) Chrysene	21.84	228	35327	1.656	ng/ul	97
21) Benzo(b)fluoranthene	23.61	252	34258	1.688	ng/ul	81
22) Benzo(k)fluoranthene	23.66	252	32127	1.730	ng/ul#	84
23) Benzo(a)pyrene	24.31	252	31122	1.717	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.23	276	35120	1.767	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.25	278	28511	1.754	ng/ul#	86
26) Benzo(g,h,i)perylene	28.10	276	29460	1.751	ng/ul#	92

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

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