

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031318\
 Data File : BE095751.D
 Acq On : 13 Mar 2018 15:48
 Operator : SJ/JU
 Sample : SST0.412
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.412

Quant Time: Mar 13 16:25:42 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	1787	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5189	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	2896	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6490	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6486	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	5899	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3532	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8249	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	5876	0.404	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	4016	0.414	ng/ul	97
7) Acenaphthylene	14.73	152	5824	0.418	ng/ul	98
8) Acenaphthene	15.07	153	4362	0.404	ng/ul	93
9) Fluorene	16.04	166	4755	0.427	ng/ul#	92
11) Pentachlorophenol	17.37	266	285	0.227	ng/ul	93
12) Phenanthrene	17.75	178	7835	0.359	ng/ul#	89
13) Anthracene	17.84	178	7319	0.366	ng/ul#	91
15) Fluoranthene	19.72	202	9503	0.336	ng/ul	84
17) Pyrene	20.07	202	9924	0.456	ng/ul#	78
18) Benzo(a)anthracene	21.77	228	8452	0.417	ng/ul#	84
19) Chrysene	21.84	228	8654	0.411	ng/ul	96
21) Benzo(b)fluoranthene	23.61	252	8169	0.393	ng/ul	85
22) Benzo(k)fluoranthene	23.66	252	7793	0.409	ng/ul#	85
23) Benzo(a)pyrene	24.31	252	7414	0.399	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.23	276	8226	0.404	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.25	278	6611	0.397	ng/ul#	88
26) Benzo(g,h,i)perylene	28.10	276	7133	0.413	ng/ul#	94

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

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