

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050318\
 Data File : BE096320.D
 Acq On : 3 May 2018 10:45
 Operator : SJ/JU
 Sample : SST0.472
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.472

Quant Time: May 03 11:38:39 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE050318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 03 11:32:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	581	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2241	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	1306	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3696	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	4855	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3956	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1515	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.63	212	5879	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	2327	0.373	ng/ul#	89
5) 2-Methylnaphthalene	12.81	142	1538	0.368	ng/ul	99
7) Acenaphthylene	14.67	152	2491	0.370	ng/ul	99
8) Acenaphthene	15.00	153	1795	0.370	ng/ul#	94
9) Fluorene	15.97	166	2492	0.388	ng/ul	88
11) Pentachlorophenol	17.31	266	159	0.179	ng/ul	97
12) Phenanthrene	17.69	178	4149	0.381	ng/ul#	88
13) Anthracene	17.78	178	3958	0.363	ng/ul#	92
15) Fluoranthene	19.66	202	5907	0.355	ng/ul	83
17) Pyrene	20.02	202	2104	0.216	ng/ul#	77
18) Benzo(a)anthracene	21.73	228	2876	0.196	ng/ul#	86
19) Chrysene	21.78	228	5462	0.404	ng/ul	95
21) Benzo(b)fluoranthene	23.54	252	5042	0.337	ng/ul	88
22) Benzo(k)fluoranthene	23.59	252	4852	0.339	ng/ul#	91
23) Benzo(a)pyrene	24.23	252	4750	0.365	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	27.13	276	5197	0.296	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.14	278	4237	0.289	ng/ul	94
26) Benzo(g,h,i)perylene	27.99	276	4446	0.291	ng/ul	97

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

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