

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031318\
 Data File : BE095749.D
 Acq On : 13 Mar 2018 14:33
 Operator : SJ/JU
 Sample : SST0.410
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.410

Quant Time: Mar 13 16:25:20 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	8.42	152	72892	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	257279	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	152238	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	333226	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	336560	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	334842	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3591	0.01	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8197	0.01	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	6058	0.008	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	4055	0.008	ng/ul	100
7) Acenaphthylene	14.74	152	5905	0.008	ng/ul	98
8) Acenaphthene	15.07	153	4737	0.008	ng/ul	93
9) Fluorene	16.04	166	4775	0.008	ng/ul	95
11) Pentachlorophenol	17.37	266	304	0.005	ng/ul	92
12) Phenanthrene	17.75	178	8483	0.008	ng/ul#	89
13) Anthracene	17.84	178	7610	0.007	ng/ul#	92
15) Fluoranthene	19.72	202	9414	0.006	ng/ul	85
17) Pyrene	20.07	202	10376	0.009	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	9766	0.009	ng/ul#	86
19) Chrysene	21.84	228	9287	0.008	ng/ul	97
21) Benzo(b)fluoranthene	23.61	252	8307	0.007	ng/ul	84
22) Benzo(k)fluoranthene	23.66	252	7739	0.007	ng/ul#	86
23) Benzo(a)pyrene	24.30	252	7393	0.007	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.23	276	8333	0.007	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.25	278	6767	0.007	ng/ul#	89
26) Benzo(g,h,i)perylene	28.10	276	7215	0.007	ng/ul#	94

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

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