

Data Path : U:\HPCHEM1\BNA E\DATA\BE011918\
 Data File : BE095074.D
 Acq On : 19 Jan 2018 14:02
 Operator : SJ/JU
 Sample : SST0.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.403

Quant Time: Jan 19 18:07:22 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE011918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 19 18:03:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4388	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	13988	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	7674	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	18455	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	17131	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	13749	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	9540	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	22987	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.34	128	15471	0.418	ng/ul#	89
5) 2-Methylnaphthalene	12.91	142	10819	0.416	ng/ul	99
7) Acenaphthylene	14.77	152	14371	0.410	ng/ul	95
8) Acenaphthene	15.11	153	11208	0.417	ng/ul	93
9) Fluorene	16.07	166	12027	0.419	ng/ul	93
11) Pentachlorophenol	17.41	266	283	0.268	ng/ul	90
12) Phenanthrene	17.78	178	19499	0.393	ng/ul#	80
13) Anthracene	17.87	178	16772	0.378	ng/ul#	85
15) Fluoranthene	19.76	202	25248	0.392	ng/ul	84
17) Pyrene	20.11	202	25276	0.409	ng/ul#	82
18) Benzo(a)anthracene	21.82	228	19527	0.405	ng/ul#	84
19) Chrysene	21.89	228	21888	0.415	ng/ul	98
21) Benzo(b)fluoranthene	23.68	252	19646	0.422	ng/ul	84
22) Benzo(k)fluoranthene	23.73	252	17772	0.405	ng/ul#	87
23) Benzo(a)pyrene	24.38	252	16370	0.405	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.33	276	14598	0.396	ng/ul#	81
25) Dibenzo(a,h)anthracene	27.36	278	10303	0.383	ng/ul	91
26) Benzo(g,h,i)perylene	28.22	276	14536	0.406	ng/ul	97

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

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