

Data Path : U:\HPCHEM1\BNA E\DATA\BE012618\
 Data File : BE095121.D
 Acq On : 26 Jan 2018 11:16
 Operator : SJ/JU
 Sample : SST00.405
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.405

Quant Time: Jan 26 12:04:30 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 26 12:00:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	3242	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	9619	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5219	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	12054	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	15761	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13947	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	6314	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	18203	0.49	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	10722	0.421	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	7158	0.403	ng/ul	100
7) Acenaphthylene	14.75	152	10020	0.429	ng/ul	96
8) Acenaphthene	15.08	153	7840	0.431	ng/ul	95
9) Fluorene	16.03	166	8073	0.416	ng/ul#	71
11) Pentachlorophenol	17.36	266	897	1.258	ng/ul	95
12) Phenanthrene	17.75	178	16737	0.523	ng/ul#	86
13) Anthracene	17.85	178	15060	0.539	ng/ul#	91
15) Fluoranthene	19.73	202	20728	0.507	ng/ul	82
17) Pyrene	20.08	202	20365	0.355	ng/ul#	80
18) Benzo(a)anthracene	21.79	228	19896	0.456	ng/ul#	84
19) Chrysene	21.85	228	20838	0.428	ng/ul	98
21) Benzo(b)fluoranthene	23.62	252	19723	0.420	ng/ul	83
22) Benzo(k)fluoranthene	23.68	252	18120	0.418	ng/ul#	84
23) Benzo(a)pyrene	24.32	252	17505	0.437	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.25	276	19311	0.543	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	15901	0.625	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	16442	0.464	ng/ul#	92

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

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