

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096851.D
 Acq On : 6 Jul 2018 12:58
 Operator : SJ/JU
 Sample : SST0.245
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.245

Quant Time: Jul 06 14:58:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 06 14:58:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2701	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	12982	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	6895	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	19955	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	15330	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	13440	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4159	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9648	0.17	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.19	128	6299	0.190	ng/ul#	92
5) 2-Methylnaphthalene	11.81	142	4091	0.190	ng/ul	98
7) Acenaphthylene	13.74	152	5490	0.183	ng/ul	98
8) Acenaphthene	14.09	153	4639	0.187	ng/ul#	91
9) Fluorene	15.08	166	4692	0.177	ng/ul	95
11) Pentachlorophenol	16.43	266	196	0.253	ng/ul#	90
12) Phenanthrene	16.82	178	10262	0.183	ng/ul#	90
13) Anthracene	16.91	178	8266	0.178	ng/ul#	92
15) Fluoranthene	18.84	202	10554	0.170	ng/ul	84
17) Pyrene	19.21	202	10367	0.193	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	6815	0.185	ng/ul#	85
19) Chrysene	21.02	228	9601	0.187	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	6936	0.175	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	7442	0.180	ng/ul#	88
23) Benzo(a)pyrene	23.11	252	5859	0.165	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.51	276	5782	0.162	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.53	278	4572	0.165	ng/ul	93
26) Benzo(g,h,i)perylene	26.20	276	5995	0.169	ng/ul	96

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

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