

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM100820\
 Data File : BM027644.D
 Acq On : 08 Oct 2020 11:53
 Operator : JU/CG
 Sample : SSTDO.102
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDO.1102

Quant Time: Oct 08 13:43:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM100820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 13:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	824	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	2278	0.40	ng/ul	0.00
7) Acenaphthene-d10	14.42	164	1509	0.40	ng/ul	0.00
11) Phenanthrene-d10	17.16	188	3753	0.40	ng/ul	0.00
17) Chrysene-d12	21.34	240	3831	0.40	ng/ul	0.00
21) Perylene-d12	23.59	264	3882	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.16	152	424	0.09	ng/ul	0.00
15) Fluoranthene-d10	19.18	212	1231	0.10	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.62	128	672	0.106	ng/ul#	84
5) 2-Methylnaphthalene	12.23	142	481	0.099	ng/ul	94
6) 1-Methylnaphthalene	12.45	142	307	0.055	ng/ul	94
8) Acenaphthylene	14.13	152	673	0.099	ng/ul#	85
9) Acenaphthene	14.48	153	527	0.104	ng/ul	94
10) Fluorene	15.46	166	663	0.099	ng/ul#	98
12) Pentachlorophenol	16.81	266	157	0.107	ng/ul	97
13) Phenanthrene	17.20	178	1143	0.102	ng/ul#	88
14) Anthracene	17.29	178	1038	0.097	ng/ul#	88
16) Fluoranthene	19.21	202	1508	0.096	ng/ul	88
18) Pyrene	19.58	202	1538	0.106	ng/ul#	88
19) Benzo(a)anthracene	21.32	228	1520	0.101	ng/ul	95
20) Chrysene	21.37	228	1626	0.105	ng/ul	94
22) Benzo(b)fluoranthene	22.91	252	1745	0.109	ng/ul#	62
23) Benzo(k)fluoranthene	22.96	252	1777	0.113	ng/ul#	65
24) Benzo(a)pyrene	23.49	252	1504	0.105	ng/ul#	56
25) Indeno(1,2,3-cd)pyrene	25.87	276	1969	0.104	ng/ul#	98
26) Dibenzo(a,h)anthracene	25.87	278	1614	0.102	ng/ul#	66
27) Benzo(a,h,i)perylene	26.55	276	1737	0.107	ng/ul#	83

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

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