

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033172.D
 Acq On : 19 Nov 2021 12:40
 Operator : CG/JU
 Sample : SST0.257
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.2009

Quant Time: Nov 19 15:05:27 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 14:17:00 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	2331	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	6238	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.580	164	3469	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	7352	0.400	ng/ul	0.00
17) Chrysene-d12	21.470	240	5851	0.400	ng/ul	0.00
23) Perylene-d12	23.813	264	4969	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.423	96	479	0.161	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.343	152	1724	0.195	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	3474	0.196	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.454	88	482	0.158	ng/ul#	76
5) Naphthalene	10.814	128	3928	0.217	ng/ul	99
7) 2-Methylnaphthalene	12.415	142	2488	0.198	ng/ul	98
8) 1-Methylnaphthalene	12.635	142	2437	0.198	ng/ul	99
10) Acenaphthylene	14.307	152	3226	0.214	ng/ul#	97
11) Acenaphthene	14.644	153	2718	0.206	ng/ul	100
12) Fluorene	15.626	166	3016	0.194	ng/ul	99
14) Pentachlorophenol	16.962	266	490	0.256	ng/ul	94
15) Phenanthrene	17.358	178	5425	0.222	ng/ul	96
16) Anthracene	17.448	178	4335	0.202	ng/ul	96
19) Fluoranthene	19.365	202	5914	0.216	ng/ul	99
20) Pyrene	19.727	202	5847	0.205	ng/ul	99
21) Benzo(a)anthracene	21.456	228	4258	0.197	ng/ul	97
22) Chrysene	21.509	228	4757	0.195	ng/ul	99
24) Benzo(b)fluoranthene	23.103	252	4575	0.201	ng/ul	78
25) Benzo(k)fluoranthene	23.151	252	4619	0.189	ng/ul#	79
26) Benzo(a)pyrene	23.711	252	3750	0.200	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	26.211	276	4856	0.225	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.226	278	3777	0.221	ng/ul#	90
29) Benzo(g,h,i)perylene	26.943	276	4335	0.231	ng/ul	96

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

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