

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121421\
 Data File : BM033469.D
 Acq On : 15 Dec 2021 09:09
 Operator : CG/JU
 Sample : PB141279BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS279

Quant Time: Dec 16 02:32:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 15 01:00:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	578	0.400	ng/ul	0.00
4) Naphthalene-d8	10.697	136	1435	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.523	164	940	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.264	188	2258	0.400	ng/ul	-0.02
17) Chrysene-d12	21.426	240	2786	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	2297	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.361	96	388	0.828	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.290	152	833	0.459	ng/ul	-0.01
18) Fluoranthene-d10	19.284	212	3130	0.516	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.393	88	1264	2.065	ng/ul#	82
5) Naphthalene	10.747	128	1979	0.377	ng/ul	99
7) 2-Methylnaphthalene	12.361	142	1218	0.407	ng/ul	92
8) 1-Methylnaphthalene	12.573	142	1235	0.400	ng/ul	96
10) Acenaphthylene	14.246	152	2025	0.385	ng/ul	97
11) Acenaphthene	14.583	153	1619	0.392	ng/ul	98
12) Fluorene	15.577	166	1792	0.406	ng/ul	99
14) Pentachlorophenol	16.927	266	454	0.560	ng/ul	98
15) Phenanthrene	17.305	178	2819	0.382	ng/ul	98
16) Anthracene	17.406	178	2527	0.388	ng/ul	100
19) Fluoranthene	19.311	202	3627	0.495	ng/ul#	94
20) Pyrene	19.673	202	3766	0.466	ng/ul#	90
21) Benzo(a)anthracene	21.414	228	2803	0.452	ng/ul	99
22) Chrysene	21.465	228	3421	0.442	ng/ul	97
24) Benzo(b)fluoranthene	23.037	252	2960	0.485	ng/ul	96
25) Benzo(k)fluoranthene	23.085	252	3355	0.479	ng/ul	98
26) Benzo(a)pyrene	23.640	252	2539	0.432	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.109	276	3120	0.475	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.131	278	2599	0.478	ng/ul	97
29) Benzo(g,h,i)perylene	26.823	276	2780	0.470	ng/ul	96

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

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