

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102521\
 Data File : BM032734.D
 Acq On : 26 Oct 2021 18:59
 Operator : CG/JU
 Sample : PB140179BS
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS179

Quant Time: Oct 27 03:03:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 26 17:09:29 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.543	152	3297	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	11747	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.194	164	5414	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.929	188	8475	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.108	240	6432	0.400	ng/ul	0.00
23) Perylene-d12	23.237	264	5887	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.916	96	3211	0.869	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.918	152	5745	0.367	ng/ul	0.00
18) Fluoranthene-d10	18.958	212	7577	0.374	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.951	88	9171	2.253	ng/ul#	80
5) Naphthalene	10.370	128	13728	0.385	ng/ul	100
7) 2-Methylnaphthalene	11.994	142	8073	0.346	ng/ul	100
8) 1-Methylnaphthalene	12.214	142	7577	0.329	ng/ul	100
10) Acenaphthylene	13.915	152	7871	0.341	ng/ul	99
11) Acenaphthene	14.255	153	7914	0.375	ng/ul	98
12) Fluorene	15.244	166	8010	0.337	ng/ul	99
14) Pentachlorophenol	16.606	266	1680	0.771	ng/ul	99
15) Phenanthrene	16.970	178	10641	0.373	ng/ul	100
16) Anthracene	17.064	178	8354	0.334	ng/ul	98
19) Fluoranthene	18.988	202	11163	0.351	ng/ul	99
20) Pyrene	19.351	202	11521	0.352	ng/ul	99
21) Benzo(a)anthracene	21.094	228	8570	0.341	ng/ul	99
22) Chrysene	21.145	228	10953	0.396	ng/ul	99
24) Benzo(b)fluoranthene	22.607	252	10796	0.397	ng/ul	86
25) Benzo(k)fluoranthene	22.648	252	12058	0.445	ng/ul#	87
26) Benzo(a)pyrene	23.147	252	8918	0.391	ng/ul#	79
27) Indeno(1,2,3-cd)pyrene	25.340	276	12164	0.436	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.347	278	9605	0.437	ng/ul#	93
29) Benzo(g,h,i)perylene	25.984	276	11515	0.474	ng/ul	98

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

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