

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071321\
 Data File : BM030963.D
 Acq On : 14 Jul 2021 16:54
 Operator : CG/JU
 Sample : PB137499BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS499

Quant Time: Jul 14 17:39:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM071321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 13 14:01:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	1027	0.400	ng/ul	0.00
4) Naphthalene-d8	10.433	136	3328	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.292	164	2058	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.041	188	4269	0.400	ng/ul	-0.01
17) Chrysene-d12	21.227	240	3369	0.400	ng/ul	0.00
23) Perylene-d12	23.418	264	3176	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.237	96	741	0.702	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.025	152	1898	0.328	ng/ul	0.00
18) Fluoranthene-d10	19.070	212	4226	0.373	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.269	88	1961	1.664	ng/ul#	81
5) Naphthalene	10.482	128	3138	0.316	ng/ul	100
7) 2-Methylnaphthalene	12.097	142	2200	0.314	ng/ul	99
8) 1-Methylnaphthalene	12.317	142	2052	0.299	ng/ul	100
10) Acenaphthylene	14.011	152	2854	0.283	ng/ul	99
11) Acenaphthene	14.356	153	2316	0.318	ng/ul	98
12) Fluorene	15.346	166	2727	0.315	ng/ul	99
14) Pentachlorophenol	16.687	266	1353	0.897	ng/ul	100
15) Phenanthrene	17.082	178	4373	0.326	ng/ul	99
16) Anthracene	17.173	178	4025	0.313	ng/ul	99
19) Fluoranthene	19.100	202	5189	0.350	ng/ul	99
20) Pyrene	19.462	202	5252	0.350	ng/ul	99
21) Benzo(a)anthracene	21.212	228	4915	0.338	ng/ul	98
22) Chrysene	21.265	228	5195	0.345	ng/ul	99
24) Benzo(b)fluoranthene	22.764	252	5352	0.348	ng/ul	98
25) Benzo(k)fluoranthene	22.808	252	5593	0.362	ng/ul	97
26) Benzo(a)pyrene	23.321	252	4437	0.332	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.596	276	6497	0.359	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.613	278	5197	0.355	ng/ul	96
29) Benzo(g,h,i)perylene	26.260	276	5478	0.357	ng/ul	98

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

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