

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120524\
 Data File : BM048888.D
 Acq On : 07 Dec 2024 03:40
 Operator : RC/JU
 Sample : PB165445BS
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS445

Quant Time: Dec 07 04:09:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM120324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 00:05:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	4802	0.400	ng/ul	0.00
4) Naphthalene-d8	10.358	136	13733	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.228	164	7714	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.975	188	15008	0.400	ng/ul	0.00
17) Chrysene-d12	21.170	240	9144	0.400	ng/ul	0.00
23) Perylene-d12	23.344	264	8530	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	4229	0.689	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.958	152	7428	0.449	ng/ul	0.00
18) Fluoranthene-d10	19.006	212	13795	0.489	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.193	88	11997	1.653	ng/ul#	75
5) Naphthalene	10.407	128	13999	0.403	ng/ul	99
7) 2-Methylnaphthalene	12.030	142	8377	0.418	ng/ul	98
8) 1-Methylnaphthalene	12.250	142	8668	0.408	ng/ul	99
10) Acenaphthylene	13.941	152	12742	0.377	ng/ul	98
11) Acenaphthene	14.288	153	9736	0.399	ng/ul	98
12) Fluorene	15.283	166	10569	0.418	ng/ul	99
14) Pentachlorophenol	16.629	266	1715	1.322	ng/ul#	4
15) Phenanthrene	17.013	178	16466	0.433	ng/ul	98
16) Anthracene	17.106	178	14069	0.415	ng/ul	97
19) Fluoranthene	19.038	202	18163	0.432	ng/ul	99
20) Pyrene	19.401	202	18344	0.419	ng/ul	99
21) Benzo(a)anthracene	21.155	228	12252	0.475	ng/ul	98
22) Chrysene	21.205	228	16012	0.392	ng/ul	99
24) Benzo(b)fluoranthene	22.698	252	12934	0.540	ng/ul	94
25) Benzo(k)fluoranthene	22.739	252	14962	0.445	ng/ul	96
26) Benzo(a)pyrene	23.250	252	11118	0.453	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.504	276	14904	0.459	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.524	278	11352	0.468	ng/ul	93
29) Benzo(g,h,i)perylene	26.155	276	12398	0.440	ng/ul	97

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

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