

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120524\
 Data File : BM048857.D
 Acq On : 06 Dec 2024 05:24
 Operator : RC/JU
 Sample : PB165399BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS399

Quant Time: Dec 06 06:14:36 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	7.598	152	4069	0.400	ng/ul	0.00
4) Naphthalene-d8	10.359	136	11126	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.229	164	5696	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.972	188	10855	0.400	ng/ul	0.00
17) Chrysene-d12	21.172	240	8317	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	9160	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.172	96	3748	0.721	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.959	152	5236	0.391	ng/ul	0.00
18) Fluoranthene-d10	19.007	212	9983	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.202	88	9894	1.608	ng/ul#	78
5) Naphthalene	10.409	128	10143	0.361	ng/ul	99
7) 2-Methylnaphthalene	12.031	142	5793	0.357	ng/ul	98
8) 1-Methylnaphthalene	12.251	142	6137	0.357	ng/ul	100
10) Acenaphthylene	13.942	152	8630	0.346	ng/ul	99
11) Acenaphthene	14.289	153	6438	0.357	ng/ul	97
12) Fluorene	15.284	166	6898	0.370	ng/ul	98
14) Pentachlorophenol	16.630	266	1087	1.158	ng/ul#	6
15) Phenanthrene	17.015	178	10739	0.390	ng/ul	100
16) Anthracene	17.108	178	9351	0.381	ng/ul	98
19) Fluoranthene	19.039	202	12968	0.339	ng/ul	99
20) Pyrene	19.402	202	13757	0.346	ng/ul	100
21) Benzo(a)anthracene	21.157	228	9790	0.417	ng/ul	98
22) Chrysene	21.207	228	13129	0.353	ng/ul	99
24) Benzo(b)fluoranthene	22.697	252	11144	0.433	ng/ul	98
25) Benzo(k)fluoranthene	22.741	252	13784	0.382	ng/ul	99
26) Benzo(a)pyrene	23.250	252	10519	0.399	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.507	276	14630	0.420	ng/ul	99
28) Dibenzo(a,h)anthracene	25.524	278	11414	0.438	ng/ul	97
29) Benzo(g,h,i)perylene	26.154	276	12126	0.401	ng/ul	97

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

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