

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122624\
 Data File : BM049224.D
 Acq On : 26 Dec 2024 13:52
 Operator : RC/JU
 Sample : PB165772BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS772

Quant Time: Dec 26 21:57:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 26 11:35:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	3600	0.400	ng/ul	0.00
4) Naphthalene-d8	10.315	136	10709	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.187	164	6560	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.939	188	14367	0.400	ng/ul	0.00
17) Chrysene-d12	21.137	240	13646	0.400	ng/ul	0.00
23) Perylene-d12	23.291	264	15620	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	2899	0.698	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.915	152	5425	0.377	ng/ul	0.00
18) Fluoranthene-d10	18.970	212	14041	0.363	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.172	88	7994	1.647	ng/ul#	77
5) Naphthalene	10.359	128	9440	0.352	ng/ul	98
7) 2-Methylnaphthalene	11.987	142	5772	0.340	ng/ul	99
8) 1-Methylnaphthalene	12.207	142	6192	0.356	ng/ul	97
10) Acenaphthylene	13.905	152	8926	0.328	ng/ul	99
11) Acenaphthene	14.248	153	6854	0.345	ng/ul	99
12) Fluorene	15.247	166	7811	0.349	ng/ul	97
14) Pentachlorophenol	16.597	266	2731	0.688	ng/ul	95
15) Phenanthrene	16.977	178	12736	0.339	ng/ul	99
16) Anthracene	17.069	178	11936	0.347	ng/ul	98
19) Fluoranthene	19.002	202	16861	0.342	ng/ul	100
20) Pyrene	19.365	202	18009	0.342	ng/ul	99
21) Benzo(a)anthracene	21.122	228	16098	0.339	ng/ul	99
22) Chrysene	21.172	228	18565	0.364	ng/ul	99
24) Benzo(b)fluoranthene	22.651	252	18102	0.363	ng/ul	98
25) Benzo(k)fluoranthene	22.691	252	20085	0.367	ng/ul	99
26) Benzo(a)pyrene	23.197	252	17103	0.357	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.427	276	23163	0.342	ng/ul	99
28) Dibenzo(a,h)anthracene	25.443	278	17902	0.340	ng/ul	98
29) Benzo(g,h,i)perylene	26.074	276	19003	0.349	ng/ul	99

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

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