

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120524\
 Data File : BM048858.D
 Acq On : 06 Dec 2024 05:59
 Operator : RC/JU
 Sample : PB165401BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS401

Quant Time: Dec 06 06:28:56 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	4257	0.400	ng/ul	0.00
4) Naphthalene-d8	10.363	136	11625	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.228	164	6176	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.975	188	11090	0.400	ng/ul	0.00
17) Chrysene-d12	21.169	240	6887	0.400	ng/ul	0.00
23) Perylene-d12	23.338	264	6977	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	4168	0.766	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.958	152	6134	0.438	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	9920	0.467	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.202	88	11623	1.806	ng/ul#	79
5) Naphthalene	10.407	128	11763	0.400	ng/ul	99
7) 2-Methylnaphthalene	12.035	142	6783	0.400	ng/ul	99
8) 1-Methylnaphthalene	12.255	142	7170	0.399	ng/ul	100
10) Acenaphthylene	13.945	152	10179	0.376	ng/ul	99
11) Acenaphthene	14.288	153	7681	0.393	ng/ul	98
12) Fluorene	15.282	166	8132	0.402	ng/ul	99
14) Pentachlorophenol	16.633	266	1117	1.165	ng/ul#	4
15) Phenanthrene	17.017	178	11854	0.422	ng/ul	100
16) Anthracene	17.110	178	10265	0.410	ng/ul	97
19) Fluoranthene	19.038	202	12968	0.410	ng/ul	99
20) Pyrene	19.400	202	13194	0.400	ng/ul	100
21) Benzo(a)anthracene	21.155	228	8522	0.438	ng/ul	99
22) Chrysene	21.207	228	12129	0.394	ng/ul	99
24) Benzo(b)fluoranthene	22.695	252	9464	0.483	ng/ul	98
25) Benzo(k)fluoranthene	22.741	252	11951	0.435	ng/ul	100
26) Benzo(a)pyrene	23.247	252	8790	0.438	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.501	276	11924	0.449	ng/ul	98
28) Dibenzo(a,h)anthracene	25.524	278	9094	0.458	ng/ul	97
29) Benzo(g,h,i)perylene	26.155	276	10147	0.441	ng/ul	97

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

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