

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103124\
 Data File : BM048412.D
 Acq On : 02 Nov 2024 01:57
 Operator : RC/JU
 Sample : PB164490BS
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS490

Quant Time: Nov 03 21:51:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	3019	0.400	ng/ul	0.03
4) Naphthalene-d8	10.541	136	11274	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.368	164	6555	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.124	188	13639	0.400	ng/ul	0.00
17) Chrysene-d12	21.295	240	11502	0.400	ng/ul	0.00
23) Perylene-d12	23.536	264	11483	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	3796	1.024	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.157	152	6151	0.429	ng/ul	0.02
18) Fluoranthene-d10	19.142	212	14331	0.491	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.198	88	9815	2.440	ng/ul#	82
5) Naphthalene	10.590	128	11343	0.385	ng/ul	98
7) 2-Methylnaphthalene	12.229	142	6376	0.364	ng/ul	99
8) 1-Methylnaphthalene	12.416	142	7372	0.395	ng/ul	100
10) Acenaphthylene	14.095	152	9772	0.386	ng/ul	98
11) Acenaphthene	14.433	153	8189	0.403	ng/ul	99
12) Fluorene	15.446	166	8680	0.384	ng/ul	99
14) Pentachlorophenol	16.787	266	2647	0.697	ng/ul	95
15) Phenanthrene	17.171	178	12748	0.376	ng/ul	99
16) Anthracene	17.285	178	11473	0.362	ng/ul	97
19) Fluoranthene	19.174	202	17883	0.451	ng/ul	99
20) Pyrene	19.532	202	19458	0.457	ng/ul	99
21) Benzo(a)anthracene	21.283	228	13411	0.389	ng/ul	99
22) Chrysene	21.330	228	22700	0.479	ng/ul	100
24) Benzo(b)fluoranthene	22.861	252	17131	0.437	ng/ul	99
25) Benzo(k)fluoranthene	22.902	252	21886	0.471	ng/ul	100
26) Benzo(a)pyrene	23.443	252	16520	0.448	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.822	276	22369	0.434	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.853	278	17335	0.431	ng/ul	100
29) Benzo(g,h,i)perylene	26.500	276	19315	0.445	ng/ul	99

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

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