

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092524\
 Data File : BM047634.D
 Acq On : 25 Sep 2024 20:45
 Operator : RC/JU
 Sample : PB163589BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS589

Quant Time: Sep 26 02:53:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 16:27:08 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	5093	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	14578	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.456	164	7404	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.196	188	15181	0.400	ng/ul	0.00
17) Chrysene-d12	21.359	240	12858	0.400	ng/ul	0.00
23) Perylene-d12	23.639	264	13855	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.273	96	4961	0.758	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.212	152	7471	0.384	ng/ul	0.00
18) Fluoranthene-d10	19.216	212	15335	0.389	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.307	88	14497	1.851	ng/ul#	81
5) Naphthalene	10.673	128	15200	0.383	ng/ul	99
7) 2-Methylnaphthalene	12.284	142	8971	0.376	ng/ul	100
8) 1-Methylnaphthalene	12.504	142	10921	0.437	ng/ul	100
10) Acenaphthylene	14.178	152	12801	0.356	ng/ul	99
11) Acenaphthene	14.521	153	9585	0.388	ng/ul	99
12) Fluorene	15.506	166	10399	0.372	ng/ul	100
14) Pentachlorophenol	16.846	266	3396	0.744	ng/ul	99
15) Phenanthrene	17.234	178	16978	0.377	ng/ul	100
16) Anthracene	17.327	178	15026	0.375	ng/ul	99
19) Fluoranthene	19.244	202	20336	0.376	ng/ul	98
20) Pyrene	19.606	202	21674	0.376	ng/ul	99
21) Benzo(a)anthracene	21.341	228	19464	0.358	ng/ul	99
22) Chrysene	21.394	228	20744	0.367	ng/ul	99
24) Benzo(b)fluoranthene	22.949	252	21952	0.364	ng/ul	99
25) Benzo(k)fluoranthene	22.993	252	21964	0.362	ng/ul	99
26) Benzo(a)pyrene	23.536	252	19166	0.403	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.960	276	27200	0.359	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.970	278	20629	0.361	ng/ul	99
29) Benzo(g,h,i)perylene	26.668	276	22197	0.365	ng/ul	99

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

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