

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042924\
 Data File : BM045626.D
 Acq On : 29 Apr 2024 19:32
 Operator : MA/JU
 Sample : P2105-02MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC7MS

Quant Time: Apr 30 01:18:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 30 01:16:44 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.559	152	918	0.400	ng/ul	0.00
4) Naphthalene-d8	10.320	136	2790	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.200	164	1297	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	2301	0.400	ng/ul	-0.01
17) Chrysene-d12	21.184	240	2034	0.400	ng/ul	0.00
23) Perylene-d12	23.370	264	2883	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.167	96	564	0.458	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.925	152	932	0.253	ng/ul	-0.01
18) Fluoranthene-d10	19.010	212	1476	0.230	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.196	88	1473	1.244	ng/ul#	86
5) Naphthalene	10.369	128	2056	0.274	ng/ul	97
7) 2-Methylnaphthalene	11.997	142	1237	0.289	ng/ul	100
8) 1-Methylnaphthalene	12.211	142	1262	0.263	ng/ul	99
10) Acenaphthylene	13.923	152	1819	0.284	ng/ul	97
11) Acenaphthene	14.265	153	1302	0.274	ng/ul	97
12) Fluorene	15.274	166	1357	0.267	ng/ul	97
14) Pentachlorophenol	16.638	266	262	0.571	ng/ul	95
15) Phenanthrene	17.014	178	2305	0.344	ng/ul	98
16) Anthracene	17.115	178	1935	0.309	ng/ul	100
19) Fluoranthene	19.043	202	2698	0.287	ng/ul	98
20) Pyrene	19.410	202	2865	0.297	ng/ul	95
21) Benzo(a)anthracene	21.173	228	2356	0.362	ng/ul	98
22) Chrysene	21.219	228	2983	0.298	ng/ul	99
24) Benzo(b)fluoranthene	22.718	252	3408	0.351	ng/ul	96
25) Benzo(k)fluoranthene	22.765	252	3280	0.268	ng/ul	97
26) Benzo(a)pyrene	23.283	252	3062	0.304	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.565	276	4379	0.300	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.575	278	3356	0.289	ng/ul	98
29) Benzo(g,h,i)perylene	26.226	276	3830	0.292	ng/ul	98

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

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