

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042224\
 Data File : BM045507.D
 Acq On : 23 Apr 2024 10:47
 Operator : MA/JU
 Sample : P2104-05
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA7

Quant Time: Apr 24 04:49:31 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 : Wed Apr 24 04:47:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.568	152	1175	0.400	ng/ul	0.00
4) Naphthalene-d8	10.331	136	3416	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.205	164	1770	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.976	188	2952	0.400	ng/ul	-0.01
17) Chrysene-d12	21.184	240	2249	0.400	ng/ul	-0.02
23) Perylene-d12	23.373	264	3296	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.167	96	3752	2.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.947	152	35	0.008	ng/ul	0.00
18) Fluoranthene-d10	19.015	212	22	0.003	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	13.927	152	2529	0.290	ng/ul	97
11) Acenaphthene	14.265	153	234	0.036	ng/ul#	87
12) Fluorene	15.274	166	1164	0.168	ng/ul#	94
15) Phenanthrene	17.018	178	1669	0.194	ng/ul	96
16) Anthracene	17.115	178	2808	0.349	ng/ul	98
19) Fluoranthene	19.047	202	2829	0.273	ng/ul	97
20) Pyrene	19.410	202	610	0.057	ng/ul#	79
21) Benzo(a)anthracene	21.169	228	1690	0.235	ng/ul#	63
22) Chrysene	21.222	228	908	0.082	ng/ul#	66
24) Benzo(b)fluoranthene	22.721	252	1590	0.143	ng/ul#	40
25) Benzo(k)fluoranthene	22.765	252	1051	0.075	ng/ul#	8
26) Benzo(a)pyrene	23.282	252	414	0.036	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.558	276	604	0.036	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.564	278	778	0.059	ng/ul#	1
29) Benzo(g,h,i)perylene	26.225	276	462	0.031	ng/ul#	10

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

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