

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046547.D
 Acq On : 12 Jul 2024 14:47
 Operator : MA/JU
 Sample : P3087-04DL 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP9DL

Quant Time: Jul 12 15:21:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 05:09:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	1948	0.400	ng/ul	0.00
4) Naphthalene-d8	10.271	136	4751	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.155	164	3007	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.905	188	6480	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	5187	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	6370	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.126	96	77	0.048	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.866	152	192	0.025	ng/ul	0.00
18) Fluoranthene-d10	18.942	212	456	0.029	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.321	128	15134	1.252	ng/ul	96
10) Acenaphthylene	13.868	152	15619	1.141	ng/ul	97
11) Acenaphthene	14.215	153	9409	1.026	ng/ul	96
12) Fluorene	15.210	166	6159	0.543	ng/ul	98
15) Phenanthrene	16.947	178	10309	0.557	ng/ul	98
16) Anthracene	17.036	178	7151	0.396	ng/ul	98
19) Fluoranthene	18.970	202	13935	0.654	ng/ul	98
20) Pyrene	19.337	202	16835	0.730	ng/ul#	97
21) Benzo(a)anthracene	21.093	228	4667	0.203	ng/ul	99
22) Chrysene	21.146	228	8556	0.385	ng/ul	98
24) Benzo(b)fluoranthene	22.618	252	10991	0.432	ng/ul	96
25) Benzo(k)fluoranthene	22.659	252	9066	0.363	ng/ul	98
26) Benzo(a)pyrene	23.159	252	11552	0.552	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.370	276	15170	0.476	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.393	278	9885	0.392	ng/ul	99
29) Benzo(g,h,i)perylene	26.014	276	10532	0.414	ng/ul	99

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

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