

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046546.D
 Acq On : 12 Jul 2024 14:11
 Operator : MA/JU
 Sample : P3102-01DL2 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP3DL2

Quant Time: Jul 12 14:39:46 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	2744	0.400	ng/ul	0.00
4) Naphthalene-d8	10.271	136	6783	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.155	164	4115	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.905	188	8585	0.400	ng/ul	0.00
17) Chrysene-d12	21.108	240	6638	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	7823	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	82	0.036	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.871	152	222	0.020	ng/ul	0.00
18) Fluoranthene-d10	18.942	212	478	0.024	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.321	128	4945	0.287	ng/ul	97
10) Acenaphthylene	13.868	152	7921	0.423	ng/ul	97
11) Acenaphthene	14.215	153	12151	0.968	ng/ul	97
12) Fluorene	15.210	166	2582	0.166	ng/ul	96
15) Phenanthrene	16.943	178	24325	0.993	ng/ul	98
16) Anthracene	17.036	178	18522	0.775	ng/ul	98
19) Fluoranthene	18.970	202	14058	0.516	ng/ul#	98
20) Pyrene	19.337	202	16924	0.573	ng/ul	99
21) Benzo(a)anthracene	21.093	228	6905	0.235	ng/ul	99
22) Chrysene	21.146	228	12098	0.426	ng/ul	99
24) Benzo(b)fluoranthene	22.616	252	16090	0.515	ng/ul	96
25) Benzo(k)fluoranthene	22.659	252	11021	0.359	ng/ul	99
26) Benzo(a)pyrene	23.156	252	1787	0.069	ng/ul#	65
27) Indeno(1,2,3-cd)pyrene	25.370	276	14013	0.358	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.390	278	11295	0.365	ng/ul	97
29) Benzo(g,h,i)perylene	26.010	276	9878	0.316	ng/ul	100

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

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