

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046555.D
 Acq On : 12 Jul 2024 21:19
 Operator : MA/JU
 Sample : P3088-01DL 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP1DL

Quant Time: Jul 13 00:36:13 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	3310	0.400	ng/ul	0.00
4) Naphthalene-d8	10.271	136	8554	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.151	164	5365	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.905	188	11286	0.400	ng/ul	0.00
17) Chrysene-d12	21.108	240	8661	0.400	ng/ul	0.00
23) Perylene-d12	23.247	264	10056	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.118	96	101	0.037	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.866	152	434	0.031	ng/ul	0.00
18) Fluoranthene-d10	18.942	212	993	0.038	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.315	128	22267	1.023	ng/ul	96
10) Acenaphthylene	13.868	152	38140	1.562	ng/ul	96
11) Acenaphthene	14.215	153	32175	1.966	ng/ul	96
12) Fluorene	15.210	166	2325	0.115	ng/ul	96
15) Phenanthrene	16.943	178	41072	1.275	ng/ul	98
16) Anthracene	17.036	178	48863	1.555	ng/ul	97
19) Fluoranthene	18.970	202	41440	1.165	ng/ul	99
20) Pyrene	19.332	202	7699	0.200	ng/ul#	97
21) Benzo(a)anthracene	21.090	228	6850	0.179	ng/ul	98
22) Chrysene	21.143	228	4421	0.119	ng/ul	97
24) Benzo(b)fluoranthene	22.613	252	15123	0.377	ng/ul	93
25) Benzo(k)fluoranthene	22.654	252	14150	0.359	ng/ul#	94
26) Benzo(a)pyrene	23.154	252	26620	0.805	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	25.366	276	41191	0.819	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.383	278	21180	0.532	ng/ul#	92
29) Benzo(g,h,i)perylene	26.004	276	8624	0.215	ng/ul#	96

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

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