

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046542.D
 Acq On : 12 Jul 2024 11:40
 Operator : MA/JU
 Sample : P3035-05DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP0DL

Quant Time: Jul 12 12:18:05 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	7.514	152	2527	0.400	ng/ul	0.00
4) Naphthalene-d8	10.270	136	6378	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.154	164	4022	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.908	188	8480	0.400	ng/ul	0.00
17) Chrysene-d12	21.114	240	6606	0.400	ng/ul	0.00
23) Perylene-d12	23.256	264	7317	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.126	96	71	0.034	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.876	152	213	0.021	ng/ul	0.00
18) Fluoranthene-d10	18.945	212	738	0.037	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.319	128	19460	1.199	ng/ul	96
10) Acenaphthylene	13.867	152	18168	0.992	ng/ul	97
11) Acenaphthene	14.219	153	14704	1.199	ng/ul	96
12) Fluorene	15.213	166	5255	0.346	ng/ul	98
15) Phenanthrene	16.946	178	14557	0.602	ng/ul	98
16) Anthracene	17.039	178	17206	0.729	ng/ul	98
19) Fluoranthene	18.973	202	30225	1.114	ng/ul	99
20) Pyrene	19.340	202	9495	0.323	ng/ul	99
21) Benzo(a)anthracene	21.097	228	17010	0.582	ng/ul	99
22) Chrysene	21.149	228	6176	0.218	ng/ul	99
24) Benzo(b)fluoranthene	22.619	252	20065	0.687	ng/ul	95
25) Benzo(k)fluoranthene	22.663	252	18124	0.632	ng/ul	95
26) Benzo(a)pyrene	23.163	252	9179	0.382	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.377	276	10902	0.298	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.397	278	2979	0.103	ng/ul#	79
29) Benzo(g,h,i)perylene	26.021	276	19442	0.666	ng/ul	98

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

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