

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046541.D
 Acq On : 12 Jul 2024 10:36
 Operator : MA/JU
 Sample : P3035-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP0

Quant Time: Jul 12 11:09:02 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	1189	0.400	ng/ul	0.00
4) Naphthalene-d8	10.270	136	2836	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.154	164	2003	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.904	188	5021	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	6348	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	8147	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	289	0.296	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.870	152	898	0.196	ng/ul	0.00
18) Fluoranthene-d10	18.940	212	4650	0.241	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.319	128	78810	10.924	ng/ul	95
10) Acenaphthylene	13.867	152	83535	9.161	ng/ul	97
11) Acenaphthene	14.214	153	68158	11.157	ng/ul	96
12) Fluorene	15.209	166	26334	3.484	ng/ul	97
15) Phenanthrene	16.946	178	77170	5.386	ng/ul	98
16) Anthracene	17.034	178	99104	7.088	ng/ul	97
19) Fluoranthene	18.973	202	204145	7.829	ng/ul#	96
20) Pyrene	19.335	202	67180	2.380	ng/ul#	97
21) Benzo(a)anthracene	21.094	228	154140	5.489	ng/ul	98
22) Chrysene	21.146	228	53402	1.965	ng/ul	98
24) Benzo(b)fluoranthene	22.619	252	200089	6.149	ng/ul	89
25) Benzo(k)fluoranthene	22.660	252	177741	5.563	ng/ul#	89
26) Benzo(a)pyrene	23.157	252	93630	3.496	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.373	276	114947	2.820	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.394	278	30761	0.953	ng/ul	94
29) Benzo(g,h,i)perylene	26.017	276	203173	6.248	ng/ul	96

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

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