

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046550.D
 Acq On : 12 Jul 2024 18:14
 Operator : MA/JU
 Sample : P3088-01DL 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP1DL

Quant Time: Jul 12 19:00:30 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	1857	0.400	ng/ul	0.00
4) Naphthalene-d8	10.270	136	4511	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.154	164	2888	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.904	188	6272	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	5341	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	6693	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	94	0.062	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.870	152	230	0.032	ng/ul	0.00
18) Fluoranthene-d10	18.940	212	560	0.034	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.319	128	11594	1.010	ng/ul	96
10) Acenaphthylene	13.867	152	20355	1.548	ng/ul	97
11) Acenaphthene	14.214	153	16992	1.929	ng/ul	98
12) Fluorene	15.209	166	1293	0.119	ng/ul	95
15) Phenanthrene	16.946	178	22424	1.253	ng/ul	99
16) Anthracene	17.034	178	26996	1.546	ng/ul	98
19) Fluoranthene	18.973	202	23465	1.070	ng/ul#	97
20) Pyrene	19.336	202	4338	0.183	ng/ul	99
21) Benzo(a)anthracene	21.094	228	4027	0.170	ng/ul	99
22) Chrysene	21.143	228	2743	0.120	ng/ul	99
24) Benzo(b)fluoranthene	22.619	252	9869	0.369	ng/ul	97
25) Benzo(k)fluoranthene	22.660	252	8983	0.342	ng/ul	100
26) Benzo(a)pyrene	23.160	252	17362	0.789	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.377	276	28541	0.852	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.394	278	14797	0.558	ng/ul	97
29) Benzo(g,h,i)perylene	26.017	276	6054	0.227	ng/ul	96

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

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