

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071224\
 Data File : BM046544.D
 Acq On : 12 Jul 2024 12:57
 Operator : MA/JU
 Sample : P3102-01DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP3DL

Quant Time: Jul 12 13:29:35 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	2756	0.400	ng/ul	0.00
4) Naphthalene-d8	10.270	136	6963	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.154	164	4368	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.904	188	9240	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	7346	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	8580	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.126	96	134	0.059	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.870	152	428	0.038	ng/ul	0.00
18) Fluoranthene-d10	18.945	212	1049	0.047	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.319	128	10351	0.584	ng/ul	97
10) Acenaphthylene	13.867	152	16875	0.849	ng/ul	98
11) Acenaphthene	14.219	153	26501	1.989	ng/ul	96
12) Fluorene	15.209	166	5642	0.342	ng/ul	97
15) Phenanthrene	16.946	178	54161	2.054	ng/ul	98
16) Anthracene	17.034	178	41818	1.625	ng/ul	98
19) Fluoranthene	18.973	202	31831	1.055	ng/ul	99
20) Pyrene	19.335	202	38639	1.183	ng/ul	99
21) Benzo(a)anthracene	21.094	228	15698	0.483	ng/ul	99
22) Chrysene	21.146	228	27048	0.860	ng/ul	99
24) Benzo(b)fluoranthene	22.616	252	35862	1.046	ng/ul	92
25) Benzo(k)fluoranthene	22.660	252	23979	0.713	ng/ul	94
26) Benzo(a)pyrene	23.160	252	3957	0.140	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	25.373	276	31317	0.730	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.390	278	24989	0.735	ng/ul	96
29) Benzo(g,h,i)perylene	26.017	276	21783	0.636	ng/ul	98

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

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