

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101623\
 Data File : BM042291.D
 Acq On : 16 Oct 2023 10:11
 Operator : MA/JU
 Sample : PB156260BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK260

Quant Time: Oct 17 01:15:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 15:38:21 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	139264	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	622061	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	381096	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	828654	20.000	ng/u1	0.00
79) Chrysene-d12	21.592	240	736920	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	796800	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	25429	5.705	ng/uL	0.00
4) Pyridine-d5	3.805	84	373095	28.264	ng/u1	0.00
7) Phenol-d5	7.163	99	426113	27.609	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	311550	29.875	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	315300	29.697	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	341810	27.965	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	162193	30.701	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	170995	30.278	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	289279	28.068	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	470388	28.988	ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	1049233	31.761	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1123702	31.079	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	179993	27.089	ng/u1	0.00
60) Fluorene-d10	15.651	176	862082	31.744	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	141566	24.388	ng/u1	0.00
73) Anthracene-d10	17.510	188	1375520	32.404	ng/u1	0.00
81) Pyrene-d10	19.798	212	1645943	36.647	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.909	264	1608733	36.783	ng/u1	0.00

Target Compounds Qvalue

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

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