

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036574.D
 Acq On : 19 Sep 2022 10:55
 Operator : CG/JU
 Sample : PB147658BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 19 22:29:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	377138	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	1659606	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1102396	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2262521	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	2022668	20.000	ng/u1	0.00
88) Perylene-d12	23.997	264	1852464	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	49899	4.960	ng/uL	0.00
4) Pyridine-d5	3.816	84	636216	21.684	ng/u1	0.00
7) Phenol-d5	7.175	99	812870	24.119	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	531324	25.959	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	654991	26.994	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	643107	25.824	ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	310958	27.682	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	294919	29.311	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	579190	24.839	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	967974	24.968	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	2069354	27.832	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2366592	26.834	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	320898	22.252	ng/u1	0.00
60) Fluorene-d10	15.639	176	1824172	27.378	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	199146	22.247	ng/u1	0.00
73) Anthracene-d10	17.486	188	2848909	27.539	ng/u1	0.00
81) Pyrene-d10	19.762	212	3132837	31.483	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	2630629	29.067	ng/u1	0.00

Target Compounds	Qvalue

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

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