

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036619.D
 Acq On : 21 Sep 2022 11:35
 Operator : CG/JU
 Sample : PB147717BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 21 22:47:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	421264	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1911666	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1212885	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2442031	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	1922313	20.000	ng/u1	0.00
88) Perylene-d12	23.991	264	1641940	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	58725	5.226	ng/uL	0.00
4) Pyridine-d5	3.816	84	534983	16.324	ng/u1	0.00
7) Phenol-d5	7.175	99	1067456	28.355	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	653123	28.568	ng/u1	0.00
11) 2-Chlorophenol-d4	7.545	132	850198	31.369	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	863055	31.026	ng/u1	0.00
21) Nitrobenzene-d5	9.186	128	429071	33.161	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	430948	37.183	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	802426	29.875	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	730941	16.368	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2551415	31.190	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	3091432	31.860	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	452164	28.498	ng/u1	0.00
60) Fluorene-d10	15.633	176	2272619	31.001	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	274752	28.437	ng/u1	0.00
73) Anthracene-d10	17.480	188	3581494	32.076	ng/u1	0.00
81) Pyrene-d10	19.762	212	3806883	40.255	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2690395	33.538	ng/u1	0.00

Target Compounds Qvalue

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

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