

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090822\
 Data File : BM036526.D
 Acq On : 08 Sep 2022 11:09
 Operator : CG/JU
 Sample : PB147480BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK480

Quant Time: Sep 08 17:48:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 02 01:33:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	182268	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	835726	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.368	164	566725	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	1128902	20.000	ng/u1	0.00
79) Chrysene-d12	21.303	240	1067717	20.000	ng/u1	0.00
88) Perylene-d12	23.597	264	888969	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	31235	6.531	ng/uL	0.00
4) Pyridine-d5	3.552	84	455700	34.552	ng/u1	0.00
7) Phenol-d5	6.898	99	484947	31.170	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	324450	33.130	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	407732	33.355	ng/u1	0.00
15) 4-Methylphenol-d8	8.445	113	404665	34.835	ng/u1	0.00
21) Nitrobenzene-d5	8.892	128	210224	33.177	ng/u1	0.00
24) 2-Nitrophenol-d4	9.610	143	228635	35.573	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.145	165	389496	33.916	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	612637	34.411	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	1312578	36.609	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	1546599	34.586	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	217271	31.982	ng/u1	0.00
60) Fluorene-d10	15.363	176	1236567	38.168	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200	182159	31.386	ng/u1	0.00
73) Anthracene-d10	17.215	188	1840212	36.660	ng/u1	0.00
81) Pyrene-d10	19.515	212	2072667	37.422	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.450	264	1622333	37.222	ng/u1	0.00

Target Compounds Qvalue

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

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