

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036618.D
 Acq On : 21 Sep 2022 11:00
 Operator : CG/JU
 Sample : PB147714BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK714

Quant Time: Sep 21 22:47:10 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	8.016	152	342290	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1564203	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1016990	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2108223	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	1748734	20.000	ng/u1	0.00
88) Perylene-d12	23.991	264	1501134	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	52773	5.780	ng/uL	0.00
4) Pyridine-d5	3.816	84	482778	18.130	ng/u1	0.00
7) Phenol-d5	7.169	99	932178	30.475	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	576253	31.021	ng/u1	0.00
11) 2-Chlorophenol-d4	7.545	132	746466	33.896	ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	755261	33.416	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	379011	35.799	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	382967	40.383	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	705809	32.115	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	652311	17.852	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2321382	33.844	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2772545	34.078	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	415972	31.267	ng/u1	0.00
60) Fluorene-d10	15.633	176	2068787	33.657	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	255884	30.678	ng/u1	0.00
73) Anthracene-d10	17.480	188	3308134	34.319	ng/u1	0.00
81) Pyrene-d10	19.762	212	3582046	41.637	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2632445	35.894	ng/u1	0.00

Target Compounds Qvalue

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

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