

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
 Data File : BP021073.D
 Acq On : 19 Jul 2024 11:24
 Operator : MA/JU
 Sample : PB162089BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 20 00:34:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 18 14:23:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	135898	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.439	136	549916	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.310	164	357031	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	799539	20.000	ng/u1	-0.01
79) Chrysene-d12	21.574	240	863950	20.000	ng/u1	0.00
88) Perylene-d12	24.886	264	1028829	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	18904	6.330	ng/uL	0.00
4) Pyridine-d5	3.652	84	251025	29.034	ng/u1	0.00
7) Phenol-d5	6.898	99	334002	30.146	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	189183	28.176	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	270332	29.614	ng/u1	0.00
15) 4-Methylphenol-d8	8.404	113	265165	28.816	ng/u1	-0.01
21) Nitrobenzene-d5	8.828	128	128963	30.193	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.539	143	150594	30.804	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.092	165	246917	27.932	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.598	131	354742	30.292	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	777103	29.941	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	908718	31.181	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	102702	27.812	ng/u1	-0.02
60) Fluorene-d10	15.316	176	692005	32.500	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.474	200	101100	20.899	ng/u1	0.00
73) Anthracene-d10	17.221	188	1151050	32.413	ng/u1	0.00
81) Pyrene-d10	19.604	212	1486364	27.828	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.656	264	1638263	32.286	ng/u1	-0.03

Target Compounds	Qvalue

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

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