

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017601.D
 Acq On : 06 Oct 2023 14:47
 Operator : MA/JU
 Sample : PB156032BL
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 23:14:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 05:55:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	93812	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	381393	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	242257	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	562819	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	578712	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	628433	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	16705	7.119	ng/uL	0.00
4) Pyridine-d5	3.722	84	232382	37.835	ng/ul	0.00
7) Phenol-d5	7.087	99	275154	34.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	168290	34.942	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	219635	35.340	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	219741	34.776	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	103524	35.282	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	117659	35.553	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	212495	34.635	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	308508	35.299	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	696464	36.856	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	738228	35.428	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	100279	30.303	ng/ul	0.00
60) Fluorene-d10	15.628	176	601563	37.531	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	84230	24.006	ng/ul	0.00
73) Anthracene-d10	17.516	188	988042	37.925	ng/ul	0.00
81) Pyrene-d10	19.769	212	1223367	38.043	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1234189	38.915	ng/ul	0.00

Target Compounds	Qvalue

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

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