

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011524\
 Data File : BP019374.D
 Acq On : 15 Jan 2024 16:42
 Operator : MA/JU
 Sample : PB158454BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK454

Quant Time: Jan 16 00:34:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	134911	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.563	136	501071	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.416	164	375904	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.181	188	926898	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	906976	20.000	ng/u1	0.00
88) Perylene-d12	23.639	264	896079	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	20052	5.620	ng/uL	-0.01
4) Pyridine-d5	3.628	84	276832	26.833	ng/u1	-0.01
7) Phenol-d5	6.958	99	315309	24.453	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.111	67	201744	26.158	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.305	132	228340	25.779	ng/u1	-0.01
15) 4-Methylphenol-d8	8.493	113	240615	23.521	ng/u1	-0.02
21) Nitrobenzene-d5	8.940	128	112460	26.678	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.657	143	134206	27.400	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.199	165	250113	25.275	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	318152	25.453	ng/u1	-0.01
46) Dimethylphthalate-d6	13.840	166	880500	26.513	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	935443	26.277	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.657	143	120429	23.615	ng/u1	-0.01
60) Fluorene-d10	15.416	176	760873	27.984	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	158640	23.182	ng/u1	0.00
73) Anthracene-d10	17.275	188	1258253	27.218	ng/u1	-0.01
81) Pyrene-d10	19.528	212	1583391	26.773	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.486	264	1345966	27.624	ng/u1	0.00

Target Compounds Qvalue

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

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