

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019351.D
 Acq On : 13 Jan 2024 03:45
 Operator : MA/JU
 Sample : PB158425BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK425

Quant Time: Jan 13 04:15:39 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	7.775	152	138056	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.569	136	541943	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.422	164	390058	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	986645	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	999159	20.000	ng/u1	0.00
88) Perylene-d12	23.639	264	1022571	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	22563	6.180	ng/uL	0.00
4) Pyridine-d5	3.634	84	323143	30.608	ng/u1	0.00
7) Phenol-d5	6.964	99	375671	28.471	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.117	67	236346	29.946	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.311	132	274705	30.307	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	305158	29.151	ng/u1	-0.01
21) Nitrobenzene-d5	8.946	128	145651	31.946	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	174893	33.014	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	324265	30.297	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	416434	30.803	ng/u1	0.00
46) Dimethylphthalate-d6	13.840	166	1166512	33.851	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	1202113	32.542	ng/u1	0.00
54) 4-Nitrophenol-d4	14.657	143	162223	30.656	ng/u1	-0.01
60) Fluorene-d10	15.416	176	939725	33.308	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	207188	28.443	ng/u1	0.00
73) Anthracene-d10	17.281	188	1580495	32.118	ng/u1	0.00
81) Pyrene-d10	19.528	212	2036852	31.263	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.486	264	1860617	33.463	ng/u1	0.00

Target Compounds	Qvalue

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

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