

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019332.D
 Acq On : 12 Jan 2024 16:23
 Operator : MA/JU
 Sample : PB158382BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK382

Quant Time: Jan 12 18:20:42 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	7.781	152	115210	20.000	ng/u1	0.00
20) Naphthalene-d8	10.575	136	465817	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	337588	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	871192	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	898418	20.000	ng/u1	0.00
88) Perylene-d12	23.639	264	930135	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	16604	5.450	ng/uL	0.00
4) Pyridine-d5	3.634	84	248688	28.227	ng/u1	0.00
7) Phenol-d5	6.969	99	281870	25.598	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	182171	27.659	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	204526	27.039	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	231423	26.491	ng/u1	-0.01
21) Nitrobenzene-d5	8.946	128	108405	27.662	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	129048	28.341	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	240719	26.167	ng/u1	0.00
31) 4-Chloroaniline-d4	10.728	131	307653	26.476	ng/u1	0.00
46) Dimethylphthalate-d6	13.845	166	827767	27.754	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	857935	26.835	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	122478	26.742	ng/u1	0.00
60) Fluorene-d10	15.422	176	712063	29.161	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	155737	24.213	ng/u1	0.00
73) Anthracene-d10	17.280	188	1242084	28.586	ng/u1	0.00
81) Pyrene-d10	19.527	212	1599303	27.300	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.486	264	1471801	29.101	ng/u1	0.00

Target Compounds Qvalue

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

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