

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103123\
 Data File : BP017835.D
 Acq On : 31 Oct 2023 17:35
 Operator : MA/JU
 Sample : PB156551BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK551

Quant Time: Nov 01 02:42:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	155348	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	579924	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	324010	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	656526	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.527	240	538113	20.000	ng/u1	0.01
88) Perylene-d12	24.092	264	554607	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	27821	6.662	ng/uL	0.02
4) Pyridine-d5	3.711	84	326319	29.847	ng/u1	0.02
7) Phenol-d5	7.052	99	408295	31.584	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	270036	33.824	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	341754	32.413	ng/u1	0.00
15) 4-Methylphenol-d8	8.610	113	315334	31.062	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	155365	33.062	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	173928	33.498	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	299906	29.446	ng/u1	0.00
31) 4-Chloroaniline-d4	10.899	131	394056	29.564	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	875626	32.193	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	966925	31.877	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	102536	26.457	ng/u1	0.01
60) Fluorene-d10	15.592	176	728249	30.973	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	107795	24.927	ng/u1	0.01
73) Anthracene-d10	17.481	188	1092746	32.571	ng/u1	0.00
81) Pyrene-d10	19.745	212	1241434	36.797	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.921	264	1048062	33.929	ng/u1	0.02

Target Compounds Qvalue

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

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