

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010923\
 Data File : BP013416.D
 Acq On : 09 Jan 2023 13:08
 Operator : CG/JU
 Sample : PB150092BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK092

Quant Time: Jan 09 14:04:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 06 01:03:32 2023
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.916	152	384263	20.000 ng/u1	0.00
20) Naphthalene-d8	10.734	136	1590882	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.563	164	1023848	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.322	188	2358889	20.000 ng/u1	0.00
79) Chrysene-d12	21.416	240	2293822	20.000 ng/u1	0.00
88) Perylene-d12	23.880	264	2528232	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.311	96	56204	5.685 ng/uL	0.00
4) Pyridine-d5	3.740	84	760728	27.589 ng/u1	0.00
7) Phenol-d5	7.081	99	914053	28.667 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	624387	31.040 ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	738883	31.073 ng/u1	0.00
15) 4-Methylphenol-d8	8.628	113	708295	29.157 ng/u1	0.00
21) Nitrobenzene-d5	9.087	128	340541	30.063 ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	379024	30.567 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	706183	28.577 ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	998978	28.695 ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	2473704	33.162 ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	2528833	30.626 ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	306906	22.478 ng/u1	0.00
60) Fluorene-d10	15.557	176	2087874	31.688 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	354863	23.725 ng/u1	0.00
73) Anthracene-d10	17.422	188	3399591	32.606 ng/u1	0.00
81) Pyrene-d10	19.657	212	4118972	33.004 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	4053155	32.637 ng/u1	0.01

Target Compounds Qvalue

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

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