

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011023\
 Data File : BP013421.D
 Acq On : 10 Jan 2023 09:56
 Operator : CG/JU
 Sample : PB150112BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK112

Quant Time: Jan 10 23:19:29 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	488270	20.000	ng/u1	0.00
20) Naphthalene-d8	10.728	136	1955555	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	1194652	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.322	188	2622942	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	2504519	20.000	ng/u1	0.00
88) Perylene-d12	23.874	264	2766511	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	59526	4.739	ng/uL	0.00
4) Pyridine-d5	3.734	84	784919	22.403	ng/u1	0.00
7) Phenol-d5	7.075	99	917023	22.634	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	644091	25.199	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	758774	25.113	ng/u1	0.00
15) 4-Methylphenol-d8	8.628	113	717099	23.231	ng/u1	0.00
21) Nitrobenzene-d5	9.087	128	339116	24.355	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	385397	25.285	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	717759	23.629	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	977508	22.843	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	2311113	26.553	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	2415464	25.071	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	278212	17.463	ng/u1	0.00
60) Fluorene-d10	15.557	176	1964472	25.552	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	321668	19.341	ng/u1	0.00
73) Anthracene-d10	17.422	188	3140728	27.090	ng/u1	0.00
81) Pyrene-d10	19.657	212	3743675	27.473	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.710	264	3714900	27.337	ng/u1	0.00

Target Compounds Qvalue

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

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