

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010423\
 Data File : BP013389.D
 Acq On : 04 Jan 2023 14:15
 Operator : CG/JU
 Sample : PB149980BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK980

Quant Time: Jan 05 04:07:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 05 04:02:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	645246	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	2627515	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	1665484	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	3596215	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	2441924	20.000	ng/u1	0.00
88) Perylene-d12	23.874	264	2278199	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	85802	5.686	ng/uL	0.00
4) Pyridine-d5	3.746	84	1191282	27.789	ng/u1	0.00
7) Phenol-d5	7.087	99	1428454	27.179	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	956806	30.179	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	1159392	28.023	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	1115472	25.903	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	546361	28.301	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	615446	28.317	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	1115280	26.419	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	1552367	26.048	ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	3494658	27.302	ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	3890929	27.664	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	455207	19.770	ng/u1	0.00
60) Fluorene-d10	15.563	176	3049146	28.158	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	493750	21.335	ng/u1	0.00
73) Anthracene-d10	17.427	188	4563058	28.125	ng/u1	0.00
81) Pyrene-d10	19.657	212	4960134	35.388	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.715	264	3236962	28.506	ng/u1	0.00

Target Compounds Qvalue

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

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