

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010323\
 Data File : BP013376.D
 Acq On : 03 Jan 2023 12:20
 Operator : CG/JU
 Sample : PB149944BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK944

Quant Time: Jan 04 01:07:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.928	152	456497	20.000 ng/u1	0.00
20) Naphthalene-d8	10.746	136	1965323	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1380959	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.334	188	3354456	20.000 ng/u1	0.00
79) Chrysene-d12	21.422	240	3373688	20.000 ng/u1	0.00
88) Perylene-d12	23.892	264	3322121	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.323	96	63112	5.911 ng/uL	0.00
4) Pyridine-d5	3.746	84	848712	27.984 ng/u1	0.00
7) Phenol-d5	7.093	99	1005905	27.053 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	673113	30.009 ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	812900	27.772 ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	816300	26.793 ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	401232	27.787 ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	452381	27.828 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	830010	26.286 ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	1196252	26.836 ng/u1	0.00
46) Dimethylphthalate-d6	13.981	166	2983456	28.111 ng/u1	0.00
49) Acenaphthylene-d8	14.269	160	3196239	27.407 ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	426729	22.352 ng/u1	0.00
60) Fluorene-d10	15.569	176	2607559	29.041 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.693	200	440563	20.409 ng/u1	0.00
73) Anthracene-d10	17.434	188	4310709	28.485 ng/u1	0.00
81) Pyrene-d10	19.669	212	5373684	27.750 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	4704154	28.409 ng/u1	0.00

Target Compounds Qvalue

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

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