

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010323\
 Data File : BP013381.D
 Acq On : 03 Jan 2023 16:57
 Operator : CG/JU
 Sample : PB149947BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK947

Quant Time: Jan 04 01:10:16 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	587138	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	2482059	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	1673336	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.328	188	3564015	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	2314145	20.000	ng/u1	-0.01
88) Perylene-d12	23.874	264	2505276	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	78807	5.739	ng/uL	0.00
4) Pyridine-d5	3.746	84	1071235	27.462	ng/u1	0.00
7) Phenol-d5	7.087	99	1285066	26.871	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	855835	29.666	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	1033087	27.441	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	1020184	26.035	ng/u1	-0.01
21) Nitrobenzene-d5	9.093	128	505991	27.746	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	570279	27.777	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	1030486	25.841	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	1470819	26.126	ng/u1	-0.01
46) Dimethylphthalate-d6	13.975	166	3502420	27.234	ng/u1	0.00
49) Acenaphthylene-d8	14.263	160	3822471	27.050	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	447993	19.366	ng/u1	-0.01
60) Fluorene-d10	15.563	176	3010422	27.670	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	472438	20.599	ng/u1	0.00
73) Anthracene-d10	17.428	188	4505791	28.023	ng/u1	0.00
81) Pyrene-d10	19.657	212	4585405	34.520	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.716	264	3476932	27.844	ng/u1	-0.02

Target Compounds Qvalue

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

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