

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017874.D
 Acq On : 02 Nov 2023 05:37
 Operator : MA/JU
 Sample : PB156564BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK564

Quant Time: Nov 02 06:54:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	166635	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.704	136	620886	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.575	164	349758	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.363	188	712980	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.510	240	584422	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	616717	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	30132	6.727	ng/uL	0.02
4) Pyridine-d5	3.705	84	363798	31.021	ng/u1	0.01
7) Phenol-d5	7.040	99	446843	32.225	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	299077	34.924	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.399	132	388183	34.323	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	357657	32.845	ng/u1	-0.01
21) Nitrobenzene-d5	9.087	128	179025	35.584	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.799	143	196839	35.410	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	335023	30.724	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.881	131	455232	31.900	ng/u1	-0.01
46) Dimethylphthalate-d6	13.992	166	992533	33.804	ng/u1	-0.01
49) Acenaphthylene-d8	14.269	160	1129699	34.501	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.828	143	112070	26.788	ng/u1	0.00
60) Fluorene-d10	15.581	176	824866	32.500	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	123747	26.350	ng/u1	0.00
73) Anthracene-d10	17.469	188	1240165	34.038	ng/u1	0.00
81) Pyrene-d10	19.727	212	1401298	38.245	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	1176108	34.240	ng/u1	0.00

Target Compounds Qvalue

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

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