

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103123\
 Data File : BP017841.D
 Acq On : 31 Oct 2023 20:59
 Operator : MA/JU
 Sample : 04950-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49D4

Quant Time: Nov 01 01:42:24 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	7.881	152	121352	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	440936	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.575	164	245498	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.369	188	500214	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	425378	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	458057	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	11836	3.628	ng/uL	0.02
4) Pyridine-d5	3.723	84	33162	3.883	ng/ul	0.03
7) Phenol-d5	7.046	99	50994	5.050	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	186773	29.948	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	173623	21.080	ng/ul	-0.01
15) 4-Methylphenol-d8	8.605	113	98932	12.475	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	108303	30.312	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.805	143	112512	28.500	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	183046	23.637	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	198937	19.630	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	647837	31.435	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	684763	29.794	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	11	0.004	ng/ul	-0.01
60) Fluorene-d10	15.581	176	531223	29.819	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	73768	22.389	ng/ul	0.00
73) Anthracene-d10	17.469	188	830153	32.476	ng/ul	0.00
81) Pyrene-d10	19.733	212	975427	36.575	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	853636	33.460	ng/ul	0.00

Target Compounds Qvalue

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

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