

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019130.D
 Acq On : 28 Dec 2023 18:14
 Operator : MA/JU
 Sample : 05986-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF7

Quant Time: Dec 28 21:49:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 28 03:15:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	125846	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	472344	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	287678	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	678097	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	781505	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	962334	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	15498	4.206	ng/uL	0.00
4) Pyridine-d5	3.658	84	100997	10.920	ng/ul	0.00
7) Phenol-d5	6.981	99	70626	6.258	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	223356	34.211	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	217562	23.913	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	137750	15.678	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	135585	33.206	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	137386	31.133	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	269775	30.275	ng/ul	0.00
31) 4-Chloroaniline-d4	10.752	131	310590	28.337	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	909979	35.812	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	964056	34.966	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	21705	5.058	ng/ul	0.00
60) Fluorene-d10	15.446	176	806463	36.738	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	133717	30.080	ng/ul	0.00
73) Anthracene-d10	17.304	188	1378156	38.917	ng/ul	0.00
81) Pyrene-d10	19.551	212	1821317	38.240	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.522	264	2216946	40.079	ng/ul	-0.01

Target Compounds Qvalue

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

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