

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012980.D
 Acq On : 08 Dec 2022 12:32
 Operator : CG/JU
 Sample : N5879-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 09 01:52:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 00:35:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	579144	20.000	ng/u1	0.00
20) Naphthalene-d8	10.798	136	2396233	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	1663767	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	3892188	20.000	ng/u1	0.00
79) Chrysene-d12	21.462	240	3633631	20.000	ng/u1	0.00
88) Perylene-d12	23.962	264	3717435	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	412	0.030	ng/uL	0.00
4) Pyridine-d5	3.787	84	417	0.011	ng/u1	0.00
7) Phenol-d5	7.128	99	55	0.001	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	129	0.005	ng/u1	0.00
11) 2-Chlorophenol-d4	7.504	132	55	0.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	30	0.001	ng/u1	-0.01
21) Nitrobenzene-d5	9.151	128	151	0.009	ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	9	0.000	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.422	165	79	0.002	ng/u1	0.01
31) 4-Chloroaniline-d4	10.928	131	179	0.003	ng/u1	0.00
46) Dimethylphthalate-d6	14.033	166	165	0.001	ng/u1	0.01
49) Acenaphthylene-d8	14.316	160	599	0.004	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	73	0.003	ng/u1	0.01
60) Fluorene-d10	15.610	176	2170	0.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	15	0.001	ng/u1	0.00
73) Anthracene-d10	17.380	188	3892188	22.166	ng/u1	-0.09
81) Pyrene-d10	19.704	212	2091	0.010	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.798	264	4332	0.023	ng/u1	0.01

Target Compounds	Qvalue

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

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