

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012058.D
 Acq On : 12 Oct 2022 01:26
 Operator : CG/JU
 Sample : N4991-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BF9

Quant Time: Oct 12 02:41:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	223193	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	901692	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	509848	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.263	188	1080356	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1176027	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	1292379	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	41045	7.251	ng/uL	0.00
4) Pyridine-d5	3.705	84	333897	22.179	ng/u1	0.00
7) Phenol-d5	7.022	99	435487	24.260	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	382060	37.497	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	477946	33.543	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	299917	21.753	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	256069	40.211	ng/u1	0.00
24) 2-Nitrophenol-d4	9.746	143	271504	42.046	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	433140	34.611	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	527528	28.605	ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	1293569	40.062	ng/u1	0.00
49) Acenaphthylene-d8	14.198	160	1574341	39.429	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	107163	16.905	ng/u1	0.00
60) Fluorene-d10	15.498	176	1193741	40.862	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	131185	22.996	ng/u1	0.00
73) Anthracene-d10	17.363	188	1866035	39.776	ng/u1	0.00
81) Pyrene-d10	19.598	212	2282823	37.890	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	2317577	36.889	ng/u1	0.00

Target Compounds Qvalue

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

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