

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012439.D
 Acq On : 01 Nov 2022 21:53
 Operator : CG/JU
 Sample : N5300-04
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA92

Quant Time: Nov 01 23:02:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	268467	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1202131	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.440	164	828199	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1833716	20.000	ng/u1	-0.01
79) Chrysene-d12	21.298	240	1832537	20.000	ng/u1	0.00
88) Perylene-d12	23.680	264	1583593	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	27360	4.103	ng/uL	0.00
4) Pyridine-d5	3.664	84	152302	7.952	ng/u1	0.00
7) Phenol-d5	6.963	99	183518	7.621	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	490041	34.091	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	508543	28.920	ng/u1	0.00
15) 4-Methylphenol-d8	8.505	113	354664	18.623	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	356874	45.779	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	389260	48.109	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	622689	35.511	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	789781	30.249	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2402743	42.380	ng/u1	0.00
49) Acenaphthylene-d8	14.134	160	2612726	40.326	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	72845	6.789	ng/u1	0.00
60) Fluorene-d10	15.439	176	2089615	43.049	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	197022	20.637	ng/u1	0.00
73) Anthracene-d10	17.304	188	3559837	45.055	ng/u1	0.00
81) Pyrene-d10	19.545	212	4210237	45.761	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	3618411	46.764	ng/u1	-0.01

Target Compounds Qvalue

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

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