

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012502.D
 Acq On : 03 Nov 2022 22:26
 Operator : CG/JU
 Sample : N5319-05
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAA9

Quant Time: Nov 04 01:03:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	329550	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1549689	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1072358	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2389206	20.000	ng/u1	-0.01
79) Chrysene-d12	21.286	240	2172828	20.000	ng/u1	-0.01
88) Perylene-d12	23.669	264	1907035	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	30967	3.784	ng/uL	0.00
4) Pyridine-d5	3.658	84	188733	8.027	ng/u1	0.00
7) Phenol-d5	6.958	99	180143	6.095	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	557862	31.615	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.311	132	548104	25.393	ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	369302	15.797	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	390044	38.813	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	418085	40.083	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	690437	30.544	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.734	131	920592	27.351	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2533324	34.510	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	2812499	33.526	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.675	143	53995	3.886	ng/u1	0.01
60) Fluorene-d10	15.428	176	2180351	34.691	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.563	200	361537	29.064	ng/u1	-0.01
73) Anthracene-d10	17.292	188	3622711	35.191	ng/u1	-0.01
81) Pyrene-d10	19.533	212	4074107	37.346	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.516	264	3330238	35.740	ng/u1	-0.01

Target Compounds Qvalue

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

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