

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012213.D
 Acq On : 20 Oct 2022 14:49
 Operator : CG/JU
 Sample : N5103-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BB2

Quant Time: Oct 20 23:23:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	410895	20.000	ng/u1	0.00
20) Naphthalene-d8	10.640	136	1667222	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.487	164	990564	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.245	188	1854185	20.000	ng/u1	0.00
79) Chrysene-d12	21.339	240	1417569	20.000	ng/u1	0.00
88) Perylene-d12	23.745	264	1465757	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	39502	3.891	ng/uL	0.00
4) Pyridine-d5	3.693	84	145878	4.986	ng/u1	0.01
7) Phenol-d5	6.999	99	168154	4.740	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	543037	25.242	ng/u1	0.00
11) 2-Chlorophenol-d4	7.363	132	524703	19.432	ng/u1	0.00
15) 4-Methylphenol-d8	8.546	113	289801	10.511	ng/u1	0.00
21) Nitrobenzene-d5	8.999	128	333054	26.988	ng/u1	0.00
24) 2-Nitrophenol-d4	9.722	143	334364	25.112	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	567565	22.725	ng/u1	0.00
31) 4-Chloroaniline-d4	10.787	131	355278	10.012	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	1926700	27.962	ng/u1	0.00
49) Acenaphthylene-d8	14.175	160	2110131	26.829	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	31450	2.436	ng/u1	0.02
60) Fluorene-d10	15.481	176	1664933	28.086	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	227391	21.557	ng/u1	0.00
73) Anthracene-d10	17.345	188	2449952	30.144	ng/u1	0.00
81) Pyrene-d10	19.586	212	2533229	31.387	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.592	264	2236541	30.427	ng/u1	0.00

Target Compounds Qvalue

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

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