

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
 Data File : BP011023.D
 Acq On : 19 Jul 2022 09:00
 Operator : CG/JU
 Sample : N3704-09
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AQ5

Quant Time: Jul 19 23:18:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	357176	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	1465864	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.339	164	820709	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	1634395	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.233	240	1621092	20.000	ng/u1	# 0.00
88) Perylene-d12	23.592	264	1648152	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	51674	5.216	ng/uL	0.00
4) Pyridine-d5	3.593	84	315033	12.302	ng/u1	0.00
7) Phenol-d5	6.857	99	220796	7.531	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	682836	36.218	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	679688	28.690	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	421076	17.940	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	410434	38.421	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	387484	36.658	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.098	165	635830	32.892	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	528209	16.575	ng/u1	0.00
46) Dimethylphthalate-d6	13.763	166	2331628	41.568	ng/u1	0.00
49) Acenaphthylene-d8	14.034	160	2712575	40.986	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	61870	5.450	ng/u1	0.00
60) Fluorene-d10	15.339	176	1977307	41.471	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	278867	34.481	ng/u1	0.00
73) Anthracene-d10	17.210	188	3281101	47.222	ng/u1	0.00
81) Pyrene-d10	19.469	212	3879815	45.570	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.439	264	3854392	48.692	ng/u1	0.01

Target Compounds Qvalue

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

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