

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011110.D
 Acq On : 26 Jul 2022 15:03
 Operator : CG/JU
 Sample : N3809-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBE69

Quant Time: Jul 26 15:55:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	451638	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	2086668	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.328	164	1265033	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	2344873	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	1466330	20.000	ng/u1	# 0.00
88) Perylene-d12	23.551	264	1283490	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	65784	5.251	ng/uL	0.00
4) Pyridine-d5	3.575	84	306005	9.450	ng/u1	0.00
7) Phenol-d5	6.840	99	345675	9.324	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	947121	39.729	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	913465	30.493	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	651377	21.947	ng/u1	0.00
21) Nitrobenzene-d5	8.828	128	602543	39.624	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	604539	40.177	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	954260	34.678	ng/u1	0.00
31) 4-Chloroaniline-d4	10.605	131	1362474	30.034	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	3749853	43.371	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	4289759	42.051	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	92974	5.313	ng/u1	0.00
60) Fluorene-d10	15.328	176	3015893	41.037	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	457638	39.441	ng/u1	0.00
73) Anthracene-d10	17.192	188	4341320	43.550	ng/u1	0.00
81) Pyrene-d10	19.451	212	4192577	54.441	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2815503	45.673	ng/u1	0.00

Target Compounds Qvalue

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

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