

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011400.D
 Acq On : 11 Aug 2022 23:07
 Operator : CG/JU
 Sample : N4128-15
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGC09

Quant Time: Aug 12 01:10:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 11 23:05:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	103075	20.000	ng/u1	0.00
20) Naphthalene-d8	10.369	136	427746	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.251	164	254569	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.028	188	552807	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.157	240	574976	20.000	ng/u1	# 0.00
88) Perylene-d12	23.463	264	641787	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11738	3.491	ng/uL	0.00
4) Pyridine-d5	3.499	84	59922	7.065	ng/u1	0.01
7) Phenol-d5	6.764	99	50349	5.520	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	160162	24.763	ng/u1	0.00
11) 2-Chlorophenol-d4	7.111	132	155641	21.746	ng/u1	0.00
15) 4-Methylphenol-d8	8.299	113	94418	13.183	ng/u1	0.00
21) Nitrobenzene-d5	8.740	128	92338	31.719	ng/u1	0.00
24) 2-Nitrophenol-d4	9.458	143	89587	33.613	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	143126	26.591	ng/u1	0.00
31) 4-Chloroaniline-d4	10.522	131	197293	20.967	ng/u1	0.00
46) Dimethylphthalate-d6	13.675	166	580866	32.886	ng/u1	0.00
49) Acenaphthylene-d8	13.940	160	657491	30.923	ng/u1	0.00
54) 4-Nitrophenol-d4	14.493	143	9772	3.029	ng/u1	0.01
60) Fluorene-d10	15.257	176	490861	32.495	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.393	200	74483	29.007	ng/u1	0.00
73) Anthracene-d10	17.128	188	868109	34.937	ng/u1	0.00
81) Pyrene-d10	19.386	212	1041130	32.867	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.316	264	1102787	34.073	ng/u1	0.00

Target Compounds	Qvalue

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

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