

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011435.D
 Acq On : 12 Aug 2022 21:33
 Operator : CG/JU
 Sample : N4165-18
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBZ0

Quant Time: Aug 12 23:18:51 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	7.569	152	79617	20.000	ng/u1	0.00
20) Naphthalene-d8	10.363	136	339061	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.245	164	192427	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.022	188	403439	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.151	240	386642	20.000	ng/u1	0.00
88) Perylene-d12	23.457	264	417415	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.075	96	9233	3.555	ng/uL	0.00
4) Pyridine-d5	3.499	84	25760	3.932	ng/u1	0.01
7) Phenol-d5	6.758	99	25948	3.683	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	112523	22.523	ng/u1	0.00
11) 2-Chlorophenol-d4	7.105	132	95989	17.363	ng/u1	0.00
15) 4-Methylphenol-d8	8.293	113	52452	9.481	ng/u1	-0.01
21) Nitrobenzene-d5	8.740	128	63144	27.364	ng/u1	0.00
24) 2-Nitrophenol-d4	9.457	143	57403	27.171	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	86564	20.289	ng/u1	0.00
31) 4-Chloroaniline-d4	10.516	131	126094	16.905	ng/u1	0.00
46) Dimethylphthalate-d6	13.669	166	386502	28.948	ng/u1	-0.01
49) Acenaphthylene-d8	13.940	160	404857	25.190	ng/u1	0.00
54) 4-Nitrophenol-d4	14.451	143	26	0.011	ng/u1	-0.03
60) Fluorene-d10	15.251	176	307128	26.898	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	44122	23.545	ng/u1	0.00
73) Anthracene-d10	17.122	188	595862	32.859	ng/u1	0.00
81) Pyrene-d10	19.386	212	675815	31.726	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.310	264	669984	31.827	ng/u1	0.00

Target Compounds Qvalue

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

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