

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011431.D
 Acq On : 12 Aug 2022 19:15
 Operator : CG/JU
 Sample : N4165-14
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBY8

Quant Time: Aug 12 23:17:17 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	75696	20.000	ng/u1	0.00
20) Naphthalene-d8	10.363	136	324290	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.251	164	189333	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.022	188	399109	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.151	240	391266	20.000	ng/u1	0.00
88) Perylene-d12	23.463	264	421840	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.075	96	11687	4.733	ng/uL	0.00
4) Pyridine-d5	3.499	84	43669	7.011	ng/u1	0.01
7) Phenol-d5	6.758	99	38332	5.722	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	119459	25.150	ng/u1	0.00
11) 2-Chlorophenol-d4	7.105	132	107753	20.501	ng/u1	0.00
15) 4-Methylphenol-d8	8.293	113	68087	12.945	ng/u1	-0.01
21) Nitrobenzene-d5	8.740	128	66112	29.955	ng/u1	0.00
24) 2-Nitrophenol-d4	9.457	143	62799	31.079	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	96831	23.730	ng/u1	0.00
31) 4-Chloroaniline-d4	10.516	131	142906	20.032	ng/u1	0.00
46) Dimethylphthalate-d6	13.675	166	399105	30.381	ng/u1	0.00
49) Acenaphthylene-d8	13.940	160	427873	27.057	ng/u1	0.00
54) 4-Nitrophenol-d4	14.487	143	5388	2.246	ng/u1	0.00
60) Fluorene-d10	15.251	176	320640	28.540	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	44647	24.084	ng/u1	0.00
73) Anthracene-d10	17.122	188	582956	32.496	ng/u1	0.00
81) Pyrene-d10	19.386	212	667515	30.966	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.310	264	656636	30.866	ng/u1	0.00

Target Compounds Qvalue

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

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