

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
 Data File : BP011195.D
 Acq On : 01 Aug 2022 12:28
 Operator : CG/JU
 Sample : PB146543BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Aug 01 22:59:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 16:05:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	282289	20.000	ng/u1	0.00
20) Naphthalene-d8	10.440	136	1220911	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.316	164	571698	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	927929	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	873438	20.000	ng/u1	# 0.00
88) Perylene-d12	23.557	264	993024	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	174	0.028	ng/uL	0.05
4) Pyridine-d5	3.558	84	392317	22.497	ng/u1	0.00
7) Phenol-d5	6.828	99	731010	29.668	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	527384	30.270	ng/u1	0.00
11) 2-Chlorophenol-d4	7.181	132	580720	29.907	ng/u1	0.00
15) 4-Methylphenol-d8	8.363	113	568382	28.200	ng/u1	0.00
21) Nitrobenzene-d5	8.810	128	291788	30.752	ng/u1	0.00
24) 2-Nitrophenol-d4	9.528	143	275512	30.313	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	450049	28.098	ng/u1	0.00
31) 4-Chloroaniline-d4	10.587	131	726118	27.964	ng/u1	0.00
46) Dimethylphthalate-d6	13.734	166	1210325	31.073	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	1553374	31.736	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	122438	17.476	ng/u1	0.00
60) Fluorene-d10	15.316	176	954743	29.170	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	87842	18.779	ng/u1	0.00
73) Anthracene-d10	17.186	188	1300638	31.811	ng/u1	0.00
81) Pyrene-d10	19.445	212	1390418	29.625	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	1550560	31.437	ng/u1	0.00

Target Compounds	Qvalue

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

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