

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010455.D
 Acq On : 21 May 2022 02:30
 Operator : CG/JU
 Sample : N2843-17
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4J7

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/23/2022
 Supervised By :mohammad ahmed 05/25/2022

Quant Time: May 21 03:50:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	166908	20.000	ng/u1	# 0.00
20) Naphthalene-d8	10.693	136	764920	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.522	164	528642	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.275	188	1158488	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.357	240	1070079	20.000	ng/u1	# 0.00
88) Perylene-d12	23.780	264	1021144	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	16073m	4.105	ng/uL	0.00
4) Pyridine-d5	3.758	84	196400	17.664	ng/u1	0.00
7) Phenol-d5	7.058	99	330220	23.572	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	233419	24.800	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.422	132	264564	24.680	ng/u1	-0.01
15) 4-Methylphenol-d8	8.599	113	281483	23.643	ng/u1	-0.01
21) Nitrobenzene-d5	9.052	128	147072	24.843	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	155368	24.765	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	271851	22.916	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.828	131	382762	21.700	ng/u1	-0.01
46) Dimethylphthalate-d6	13.934	166	935195	24.011	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	1129630	25.144	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	106282	15.331	ng/u1	0.00
60) Fluorene-d10	15.516	176	846795	24.991	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	95801	14.140	ng/u1	0.00
73) Anthracene-d10	17.375	188	1303648	24.851	ng/u1	0.00
81) Pyrene-d10	19.604	212	1549616	27.268	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1270302	24.774	ng/u1	-0.01

Target Compounds Qvalue

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

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