

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011222\
 Data File : BN018391.D
 Acq On : 12 Jan 2022 14:01
 Operator : CG/JU
 Sample : PB141977BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK977

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 01/12/2022
 Supervised By :mohammad ahmed 01/18/2022

Quant Time: Jan 12 15:05:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 07 01:31:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	294892	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.363	136	1323036	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.234	164	878560	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.975	188	1782451	20.000	ng/u1	0.00
79) Chrysene-d12	21.175	240	1855540	20.000	ng/u1	# 0.00
88) Perylene-d12	23.398	264	1928584	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.052	96	42280	5.763	ng/uL	-0.01
4) Pyridine-d5	3.458	84	620947	27.047	ng/u1	0.00
7) Phenol-d5	6.764	99	794135	27.709	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.922	67	532021	29.141	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.117	132	628414	29.934	ng/u1	-0.01
15) 4-Methylphenol-d8	8.299	113	626014	27.673	ng/u1	-0.02
21) Nitrobenzene-d5	8.728	128	280228	28.438	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.452	143	308091	29.250	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.993	165	531572m	27.777	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.505	131	841535	28.613	ng/u1	-0.01
46) Dimethylphthalate-d6	13.652	166	2021701	31.682	ng/u1	0.00
49) Acenaphthylene-d8	13.922	160	2454912	30.432	ng/u1	0.00
54) 4-Nitrophenol-d4	14.451	143	246294	20.226	ng/u1	0.00
60) Fluorene-d10	15.228	176	1697994	31.538	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	236537	23.452	ng/u1	0.00
73) Anthracene-d10	17.075	188	2752001	33.336	ng/u1	0.00
81) Pyrene-d10	19.375	212	3182190	29.896	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.257	264	3253002	33.892	ng/u1	0.00

Target Compounds Qvalue

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

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