

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011922\
 Data File : BN018417.D
 Acq On : 19 Jan 2022 15:04
 Operator : CG/JU
 Sample : N1129-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBLZ9

Manual Integrations
 APPROVED

Reviewed By :Jagrut
 Upadhyay
 01/20/2022

Supervised By :mohammad
 ahmed
 01/21/2022

Quant Time: Jan 20 02:34:38 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	374514	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.363	136	1677723	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.234	164	1115641	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.975	188	2253662	20.000	ng/u1	0.00
79) Chrysene-d12	21.174	240	2275319	20.000	ng/u1	# 0.00
88) Perylene-d12	23.398	264	2317049	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.058	96	25572	2.745	ng/uL	0.00
4) Pyridine-d5	3.464	84	192981	6.619	ng/u1	0.00
7) Phenol-d5	6.763	99	532093	14.619	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.922	67	365122	15.748	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.116	132	438335	16.441	ng/u1	-0.01
15) 4-Methylphenol-d8	8.299	113	369037	12.845	ng/u1	-0.02
21) Nitrobenzene-d5	8.728	128	202744	16.225	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.451	143	225200	16.860	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.993	165	388667	16.016	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.504	131	489201	13.117	ng/u1	-0.01
46) Dimethylphthalate-d6	13.645	166	1413086	17.438	ng/u1	-0.01
49) Acenaphthylene-d8	13.922	160	1728128	16.870	ng/u1	0.00
54) 4-Nitrophenol-d4	14.451	143	122849m	7.945	ng/u1	0.00
60) Fluorene-d10	15.228	176	1232204	18.023	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.351	200	139527	10.941	ng/u1	-0.01
73) Anthracene-d10	17.075	188	1888467	18.093	ng/u1	0.00
81) Pyrene-d10	19.374	212	2283037	17.491	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.257	264	2217802	19.232	ng/u1	0.00

Target Compounds Qvalue

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

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