Quantitation Report (LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030822\
Data File : BN018876.D
Acq On : 08 Mar 2022 18:19
Operator : CG/JU
Sample : N1733-27
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBRU5

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 03/09/2022
Supervised By :Yogesh Patel 03/10/2022

Quant Time: Mar 09 01:56:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN030822.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 08 14:15:20 2022
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	112889	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	532190	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	377715	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	846148	20.000	ng/ul	0.00
79) Chrysene-d12	21.457	240	943972	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	1040335	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	17685m	6.266	ng/uL	0.00
4) Pyridine-d5	3.693	84	245602	28.437	ng/ul	0.00
7) Phenol-d5	7.058	99	387223	35.134	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	253934	38.252	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	287798	36.437	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	348037	38.836	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	162700	39.800	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	183698	40.900	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	339870	39.560	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	529590	40.937	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1318796	46.115	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	1614895	45.366	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	236600	40.853	ng/ul	0.00
60) Fluorene-d10	15.528	176	1197371	48.168	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	192328	36.677	ng/ul	0.00
73) Anthracene-d10	17.375	188	1936009	47.836	ng/ul	0.00
81) Pyrene-d10	19.669	212	2362844	46.254	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	2601451	47.980	ng/ul	0.00

Target Compounds	Qvalue

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

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