

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
 Data File : BP010277.D
 Acq On : 10 May 2022 22:28
 Operator : CG/JU
 Sample : SP5897
 Misc : SFAM SURROGATE
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP5897

Quant Time: May 11 01:19:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 10 15:59:46 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
 Supervised By :mohammad ahmed 05/12/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	123719	20.000	ng/u1	0.00
20) Naphthalene-d8	10.710	136	540396	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.540	164	379070	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.292	188	859904	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.375	240	870591	20.000	ng/u1	#-0.01
88) Perylene-d12	23.810	264	906846	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	44012m	15.113	ng/uL	0.00
4) Pyridine-d5	3.764	84	636798	78.576	ng/u1	0.00
7) Phenol-d5	7.075	99	748422	65.132	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	515784	72.558	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	595898	70.414	ng/u1	0.00
15) 4-Methylphenol-d8	8.616	113	647214	68.510	ng/u1	0.00
21) Nitrobenzene-d5	9.069	128	325248	75.063	ng/u1	0.00
24) 2-Nitrophenol-d4	9.793	143	363633	73.605	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	599875	66.712	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.846	131	914045	68.855	ng/u1	-0.01
46) Dimethylphthalate-d6	13.951	166	2233418	73.956	ng/u1	0.00
49) Acenaphthylene-d8	14.234	160	2516950	69.905	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	371984	64.216	ng/u1	-0.01
60) Fluorene-d10	15.534	176	1858237	72.896	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.651	200	385653	65.492	ng/u1	-0.01
73) Anthracene-d10	17.392	188	2976030	70.673	ng/u1	-0.01
81) Pyrene-d10	19.622	212	3578113	74.374	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.657	264	3585809	74.145	ng/u1	-0.01

Target Compounds Qvalue

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

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