

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009850.D
 Acq On : 15 Apr 2022 04:47
 Operator : CG/JU
 Sample : SP5861
 Misc : SFAM SURROGATE
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP5861

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/22/2022

Quant Time: Apr 15 05:35:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	122522	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.769	136	546903	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.592	164	376255	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.339	188	837266	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.416	240	921035	20.000	ng/u1	#-0.02
88) Perylene-d12	23.880	264	908294	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	43977m	16.019	ng/uL	0.00
4) Pyridine-d5	3.793	84	602712	79.627	ng/u1	-0.01
7) Phenol-d5	7.116	99	765986	71.210	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	506464	79.088	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	622313	76.697	ng/u1	-0.01
15) 4-Methylphenol-d8	8.663	113	671221	74.750	ng/u1	0.00
21) Nitrobenzene-d5	9.116	128	340358	86.300	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.846	143	379038	87.662	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	641785	70.931	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.899	131	973286	75.525	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	2277278	77.051	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	2596541	74.156	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	398918	74.425	ng/u1	0.00
60) Fluorene-d10	15.586	176	1938454	77.965	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	358432	72.265	ng/u1	0.00
73) Anthracene-d10	17.439	188	3094780	77.200	ng/u1	-0.01
81) Pyrene-d10	19.669	212	3754388	74.636	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.727	264	3740789	79.085	ng/u1	-0.01

Target Compounds Qvalue

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

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