

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016846.D
 Acq On : 29 Aug 2023 23:02
 Operator : MA/JU
 Sample : 04148-06
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJW3

Quant Time: Aug 29 23:39:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 29 01:12:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	435156	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	1934416	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.681	164	1204727	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	2604884	20.000	ng/u1	-0.01
79) Chrysene-d12	21.539	240	2042234	20.000	ng/u1	-0.01
88) Perylene-d12	24.110	264	1832992	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	46568	4.423	ng/uL	0.00
4) Pyridine-d5	3.816	84	287082	8.922	ng/u1	0.00
7) Phenol-d5	7.169	99	268352	6.935	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	732874	29.717	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	700932	24.093	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	487226	15.425	ng/u1	0.00
21) Nitrobenzene-d5	9.222	128	446529	30.530	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	449039	29.126	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	813252	28.295	ng/u1	0.00
31) 4-Chloroaniline-d4	11.010	131	514799	11.680	ng/u1	0.00
46) Dimethylphthalate-d6	14.092	166	2810783	30.938	ng/u1	0.00
49) Acenaphthylene-d8	14.381	160	3204224	31.198	ng/u1	0.00
54) 4-Nitrophenol-d4	14.892	143	69252	3.688	ng/u1	0.00
60) Fluorene-d10	15.675	176	2479413	32.407	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	329299	21.227	ng/u1	0.00
73) Anthracene-d10	17.545	188	3930527	34.041	ng/u1	-0.01
81) Pyrene-d10	19.780	212	4337896	38.675	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.939	264	3011873	32.803	ng/u1	-0.01

Target Compounds Qvalue

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

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