

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090723\
 Data File : BP016998.D
 Acq On : 07 Sep 2023 21:03
 Operator : MA/JU
 Sample : 04123-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH931

Quant Time: Sep 08 00:41:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 06 01:52:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	352073	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1558646	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	964810	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	2029537	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	1555404	20.000	ng/u1	0.00
88) Perylene-d12	24.039	264	1508163	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	55691	5.983	ng/uL	0.00
4) Pyridine-d5	3.764	84	320181	12.039	ng/u1	0.00
7) Phenol-d5	7.128	99	267942	8.330	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.316	67	775728	38.556	ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	696585	29.923	ng/u1	0.00
15) 4-Methylphenol-d8	8.687	113	501380	20.461	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	456611	37.761	ng/u1	0.00
24) 2-Nitrophenol-d4	9.893	143	485991	37.076	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	815523	33.700	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1079065	30.906	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2996487	41.723	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	3266635	39.413	ng/u1	0.00
54) 4-Nitrophenol-d4	14.845	143	95338	6.796	ng/u1	-0.01
60) Fluorene-d10	15.634	176	2536033	41.243	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	408927	33.481	ng/u1	0.00
73) Anthracene-d10	17.504	188	3981928	42.722	ng/u1	0.00
81) Pyrene-d10	19.739	212	4321265	45.612	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.874	264	3430813	44.840	ng/u1	0.00

Target Compounds Qvalue

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

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