

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017051.D
 Acq On : 09 Sep 2023 14:34
 Operator : MA/JU
 Sample : 04227-15
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA06

Quant Time: Sep 09 23:06:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 09 01:41:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	325695	20.000	ng/u1	0.00
20) Naphthalene-d8	10.799	136	1402600	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	823498	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	1747457	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	1349542	20.000	ng/u1	0.00
88) Perylene-d12	24.027	264	1343987	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	49917	5.797	ng/uL	0.00
4) Pyridine-d5	3.764	84	405670	16.489	ng/u1	0.00
7) Phenol-d5	7.116	99	286924	9.643	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	865400	46.497	ng/u1	0.00
11) 2-Chlorophenol-d4	7.493	132	766314	35.584	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	522749	23.060	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	491240	45.145	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	492459	41.749	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	843427	38.730	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1111016	35.361	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2932289	47.836	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	3327207	47.032	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	80664	6.737	ng/u1	0.00
60) Fluorene-d10	15.622	176	2498066	47.597	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	370360	35.218	ng/u1	-0.01
73) Anthracene-d10	17.492	188	3906383	48.677	ng/u1	0.00
81) Pyrene-d10	19.733	212	4290264	52.193	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	3355006	49.206	ng/u1	0.00

Target Compounds Qvalue

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

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