

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017313.D
 Acq On : 25 Sep 2023 10:37
 Operator : MA/JU
 Sample : 04429-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKK1

Quant Time: Sep 26 01:07:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	156671	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.751	136	636900	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.592	164	371143	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	834903	20.000	ng/u1	0.00
79) Chrysene-d12	21.463	240	818521	20.000	ng/u1	-0.01
88) Perylene-d12	23.980	264	912781	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	13327	3.494	ng/uL	0.00
4) Pyridine-d5	3.740	84	90354	8.473	ng/u1	0.00
7) Phenol-d5	7.087	99	94515	7.345	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	109849	14.146	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.452	132	124013	12.152	ng/u1	-0.01
15) 4-Methylphenol-d8	8.646	113	96684	9.665	ng/u1	-0.01
21) Nitrobenzene-d5	9.128	128	63883	13.874	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.840	143	64468	12.895	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.369	165	125942	13.333	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.916	131	142596	10.422	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	393825	14.812	ng/u1	-0.01
49) Acenaphthylene-d8	14.286	160	431126	14.099	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.828	143	23114	4.990	ng/u1	0.00
60) Fluorene-d10	15.586	176	347988	15.203	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.716	200	31068	6.602	ng/u1	-0.01
73) Anthracene-d10	17.463	188	571131	15.424	ng/u1	0.00
81) Pyrene-d10	19.704	212	695340	15.709	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.815	264	682414	15.528	ng/u1	-0.02

Target Compounds Qvalue

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

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