

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017459.D
 Acq On : 29 Sep 2023 22:59
 Operator : MA/JU
 Sample : 04497-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ71

Quant Time: Sep 30 00:39:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 29 22:31:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	216324	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.787	136	884265	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.646	164	531620	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.434	188	1204607	20.000	ng/ul	-0.01
79) Chrysene-d12	21.557	240	1209880	20.000	ng/ul	-0.02
88) Perylene-d12	24.133	264	1383662	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	8363	1.588	ng/uL	0.00
4) Pyridine-d5	3.740	84	30778	2.090	ng/ul	0.00
7) Phenol-d5	7.105	99	95158	5.356	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	258985	24.154	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.469	132	273869	19.435	ng/ul	-0.01
15) 4-Methylphenol-d8	8.669	113	145162	10.509	ng/ul	-0.02
21) Nitrobenzene-d5	9.158	128	175599	27.469	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.875	143	180511	26.005	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.405	165	303983	23.180	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.957	131	178812	9.413	ng/ul	-0.01
46) Dimethylphthalate-d6	14.057	166	1097563	28.820	ng/ul	-0.01
49) Acenaphthylene-d8	14.340	160	1180497	26.951	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.887	143	26697	4.024	ng/ul	0.00
60) Fluorene-d10	15.645	176	980132	29.894	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.787	200	133874	19.718	ng/ul	-0.02
73) Anthracene-d10	17.534	188	1612311	30.179	ng/ul	-0.01
81) Pyrene-d10	19.781	212	2002681	30.609	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.963	264	1909165	28.658	ng/ul	-0.02

Target Compounds Qvalue

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

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