

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017099.D
 Acq On : 11 Sep 2023 11:57
 Operator : MA/JU
 Sample : 04262-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJD6

Quant Time: Sep 12 02:27:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 06:13:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	304560	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	1346025	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164	818205	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.387	188	1770277	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1444057	20.000	ng/ul	0.00
88) Perylene-d12	24.016	264	1410222	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	46219	5.792	ng/uL	0.00
4) Pyridine-d5	3.758	84	221829	9.219	ng/ul	0.00
7) Phenol-d5	7.111	99	192868	6.967	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	596580	32.130	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	496396	24.808	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	352887	16.113	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	318558	33.242	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	317278	32.079	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	555216	29.605	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	778629	25.891	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	2188656	37.593	ng/ul	0.00
49) Acenaphthylene-d8	14.316	160	2323201	35.044	ng/ul	0.00
54) 4-Nitrophenol-d4	14.840	143	63926	5.322	ng/ul	0.00
60) Fluorene-d10	15.616	176	1870547	37.862	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	282185	28.246	ng/ul	0.00
73) Anthracene-d10	17.487	188	3023433	39.845	ng/ul	0.00
81) Pyrene-d10	19.728	212	3422099	45.000	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	2731142	40.342	ng/ul	0.00

Target Compounds Qvalue

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

