

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017097.D
 Acq On : 11 Sep 2023 10:48
 Operator : MA/JU
 Sample : 04262-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJD4

Quant Time: Sep 12 02:27:04 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	337706	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	1474643	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	885448	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.387	188	1882389	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	1468866	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	1401144	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	45334	5.124	ng/uL	0.00
4) Pyridine-d5	3.752	84	278691	10.445	ng/u1	0.00
7) Phenol-d5	7.105	99	248754	8.103	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	680919	33.073	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	597982	26.952	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	427822	17.618	ng/u1	0.00
21) Nitrobenzene-d5	9.158	128	363568	34.630	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	374990	34.607	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	620065	30.179	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	810919	24.613	ng/u1	0.00
46) Dimethylphthalate-d6	14.034	166	2311260	36.684	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	2491640	34.731	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	78826	6.064	ng/u1	0.00
60) Fluorene-d10	15.616	176	1957792	36.618	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	307349	28.933	ng/u1	0.00
73) Anthracene-d10	17.487	188	3148244	39.019	ng/u1	0.00
81) Pyrene-d10	19.728	212	3552449	45.925	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	2733900	40.644	ng/u1	0.00

Target Compounds	Qvalue

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

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