

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017304.D
 Acq On : 23 Sep 2023 01:09
 Operator : MA/JU
 Sample : 04410-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBK36

Quant Time: Sep 23 02:07:56 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	7.934	152	157960	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	642805	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	387362	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	880458	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	829883	20.000	ng/ul	-0.01
88) Perylene-d12	23.980	264	922794	20.000	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	16453	4.278	ng/uL	0.00
4) Pyridine-d5	3.740	84	134978	12.555	ng/ul	0.00
7) Phenol-d5	7.087	99	85482	6.589	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.275	67	251796	32.161	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	260310	25.299	ng/ul	-0.01
15) 4-Methylphenol-d8	8.646	113	161513	16.013	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	158839	34.180	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	163741	32.450	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	285911	29.991	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.922	131	368710	26.700	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1043764	37.614	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	1104079	34.594	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	24457	5.059	ng/ul	0.00
60) Fluorene-d10	15.592	176	911806	38.167	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	129935	26.184	ng/ul	0.00
73) Anthracene-d10	17.463	188	1570552	40.220	ng/ul	0.00
81) Pyrene-d10	19.704	212	1932223	43.055	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.815	264	1874626	42.193	ng/ul	-0.02

Target Compounds	Qvalue

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

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