

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017170.D
 Acq On : 14 Sep 2023 11:07
 Operator : MA/JU
 Sample : 04274-01
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RT2683

Quant Time: Sep 14 22:45:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	177458	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	708711	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	400852	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	943073	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	929614	20.000	ng/u1	0.00
88) Perylene-d12	24.010	264	1058403	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	16469	3.900	ng/uL	0.00
4) Pyridine-d5	3.746	84	257	0.021	ng/u1	0.00
7) Phenol-d5	7.093	99	11	0.001	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	76	0.008	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	75	0.007	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	13	0.001	ng/u1	-0.01
21) Nitrobenzene-d5	9.140	128	79	0.016	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.857	143	12	0.002	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.222	165	35	0.003	ng/u1	-0.17
31) 4-Chloroaniline-d4	10.922	131	25	0.002	ng/u1	-0.02
46) Dimethylphthalate-d6	14.028	166	368	0.013	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	578	0.018	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	25	0.005	ng/u1	0.00
60) Fluorene-d10	15.610	176	1013	0.041	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	27	0.005	ng/u1	0.01
73) Anthracene-d10	17.381	188	943073	23.536	ng/u1	-0.11
81) Pyrene-d10	19.716	212	2410	0.049	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.839	264	2605	0.052	ng/u1	-0.01

Target Compounds Qvalue

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

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