

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014232.D
 Acq On : 22 Mar 2023 23:54
 Operator : CG/JU
 Sample : 01977-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYG95

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/23/2023
 Supervised By : Jagrut Upadhyay 03/23/2023

Quant Time: Mar 23 02:36:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 23 02:31:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	142064	20.000	ng/u1	0.00
20) Naphthalene-d8	10.863	136	631999	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.686	164	474554	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.439	188	1152005	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.521	240	1199800	20.000	ng/u1	#-0.01
88) Perylene-d12	24.045	264	1253073	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	17770m	4.726	ng/uL	0.00
4) Pyridine-d5	3.840	84	226033	21.712	ng/u1	0.00
7) Phenol-d5	7.199	99	315494	25.166	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	209826	28.698	ng/u1	0.00
11) 2-Chlorophenol-d4	7.569	132	240482	25.177	ng/u1	0.00
15) 4-Methylphenol-d8	8.752	113	261062	25.493	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	128065	25.396	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	149943	24.840	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	257052	22.632	ng/u1	0.00
31) 4-Chloroaniline-d4	11.004	131	355752	23.668	ng/u1	0.00
46) Dimethylphthalate-d6	14.086	166	1042241	25.879	ng/u1	0.00
49) Acenaphthylene-d8	14.381	160	1098548	25.749	ng/u1	0.00
54) 4-Nitrophenol-d4	14.892	143	140444m	19.788	ng/u1	0.00
60) Fluorene-d10	15.681	176	893908	26.189	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	171407	19.686	ng/u1	0.00
73) Anthracene-d10	17.539	188	1476536	26.402	ng/u1	0.00
81) Pyrene-d10	19.769	212	1803813	27.259	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.880	264	1753580	26.309	ng/u1	-0.01

Target Compounds Qvalue

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

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