

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051823\
 Data File : BP051823.D
 Acq On : 18 May 2023 16:41
 Operator : CG/JU
 Sample : 02612-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREA3

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/19/2023
 Supervised By : Jagrut Upadhyay 05/19/2023

Quant Time: May 19 01:20:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 19 01:07:51 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	312151	20.000	ng/u1	0.00
20) Naphthalene-d8	10.957	136	1436885	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.769	164	922111	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.516	188	2003050	20.000	ng/u1	0.00
79) Chrysene-d12	21.592	240	1661188	20.000	ng/u1	0.00
88) Perylene-d12	24.157	264	1592628	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	23364	2.908	ng/uL	0.00
4) Pyridine-d5	3.905	84	22318m	0.980	ng/u1	0.04
7) Phenol-d5	7.275	99	653626	22.552	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	413939	24.729	ng/u1	0.00
11) 2-Chlorophenol-d4	7.652	132	495173	23.237	ng/u1	0.00
15) 4-Methylphenol-d8	8.834	113	536551	23.233	ng/u1	0.00
21) Nitrobenzene-d5	9.305	128	263261	23.316	ng/u1	0.00
24) 2-Nitrophenol-d4	10.034	143	285113	22.694	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	512743	22.618	ng/u1	0.00
31) 4-Chloroaniline-d4	11.093	131	696195	20.756	ng/u1	0.00
46) Dimethylphthalate-d6	14.169	166	1968974	28.257	ng/u1	0.00
49) Acenaphthylene-d8	14.463	160	2142362	26.883	ng/u1	0.00
54) 4-Nitrophenol-d4	14.951	143	268726	19.861	ng/u1	0.00
60) Fluorene-d10	15.757	176	1655469	28.156	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	222850	17.654	ng/u1	0.00
73) Anthracene-d10	17.616	188	2615945	28.521	ng/u1	0.00
81) Pyrene-d10	19.839	212	3128982	32.477	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.986	264	2485601	31.440	ng/u1	0.00
Target Compounds						
16) Acetophenone	8.963	105	159564	4.427	ng/u1	93

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

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