

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016112.D
 Acq On : 09 Jul 2023 02:39
 Operator : MA/JU
 Sample : 03350-06
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW29

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/10/2023
 Supervised By :mohammad ahmed 07/10/2023

Quant Time: Jul 09 03:09:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	287175	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.852	136	1329390	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.669	164	930345	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.434	188	2145576	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.539	240	2056590	20.000	ng/u1	# 0.00
88) Perylene-d12	24.069	264	2051298	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	28816	3.610	ng/uL	0.00
4) Pyridine-d5	3.799	84	332721	16.536	ng/u1	0.00
7) Phenol-d5	7.228	99	422500	17.195	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.352	67	331009m	21.775	ng/u1	0.00
11) 2-Chlorophenol-d4	7.558	132	371855	20.048	ng/u1	0.00
15) 4-Methylphenol-d8	8.781	113	379370	19.544	ng/u1	0.01
21) Nitrobenzene-d5	9.234	128	170454	16.778	ng/u1	0.02
24) 2-Nitrophenol-d4	9.969	143	197781	16.680	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.557	165	280774	13.288	ng/u1	0.05
31) 4-Chloroaniline-d4	11.057	131	435907	16.059	ng/u1	0.04
46) Dimethylphthalate-d6	14.099	166	1327227	18.457	ng/u1	0.01
49) Acenaphthylene-d8	14.369	160	1438272	17.793	ng/u1	0.00
54) 4-Nitrophenol-d4	15.051	143	114931m	12.112	ng/u1	0.09
60) Fluorene-d10	15.681	176	1124520	18.992	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.887	200	114509m	8.678	ng/u1	0.04
73) Anthracene-d10	17.551	188	1735161	17.705	ng/u1	0.00
81) Pyrene-d10	19.775	212	2149803	16.009	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	1793026	17.328	ng/u1	-0.01

Target Compounds Qvalue

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

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