

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
 Data File : BP016177.D
 Acq On : 11 Jul 2023 03:09
 Operator : MA/JU
 Sample : 03404-18
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW54

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/11/2023
 Supervised By :mohammad ahmed 07/11/2023

Quant Time: Jul 11 03:52:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 10 23:19:51 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.017	152	191546m	20.000	ng/u1	0.00
20) Naphthalene-d8	10.846	136	893657	20.000	ng/u1	# 0.01
38) Acenaphthene-d10	14.669	164	573758	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.434	188	1266408	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.551	240	1164293	20.000	ng/u1	# 0.00
88) Perylene-d12	24.074	264	1201427	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	24667	4.633	ng/uL	0.00
4) Pyridine-d5	3.799	84	209487	15.610	ng/u1	0.00
7) Phenol-d5	7.222	99	434599	26.518	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.352	67	318436m	31.406	ng/u1	0.00
11) 2-Chlorophenol-d4	7.558	132	349134	28.221	ng/u1	0.01
15) 4-Methylphenol-d8	8.781	113	387264	29.911	ng/u1	0.02
21) Nitrobenzene-d5	9.228	128	177411	25.977	ng/u1	0.02
24) 2-Nitrophenol-d4	9.969	143	206995	25.968	ng/u1	0.03
28) 2,4-Dichlorophenol-d3	10.569	165	344806m	24.275	ng/u1	0.06
31) 4-Chloroaniline-d4	11.075	131	347506m	19.045	ng/u1	0.05
46) Dimethylphthalate-d6	14.087	166	1518272	34.236	ng/u1	0.00
49) Acenaphthylene-d8	14.363	160	1636431	32.826	ng/u1	0.00
54) 4-Nitrophenol-d4	15.087	143	128714m	21.994	ng/u1	0.09
60) Fluorene-d10	15.675	176	1317304	36.075	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.881	200	132151	16.968	ng/u1	0.02
73) Anthracene-d10	17.540	188	2075117	35.872	ng/u1	0.00
81) Pyrene-d10	19.769	212	2551351	33.559	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.898	264	2097754	34.613	ng/u1	-0.02

Target Compounds Qvalue

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

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