

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016129.D
 Acq On : 09 Jul 2023 12:50
 Operator : MA/JU
 Sample : 03356-07
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW36

Manual Integrations
 APPROVED

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

Quant Time: Jul 09 23:47:54 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	370340	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.846	136	1554583	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.669	164	940384	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.434	188	1907369	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.545	240	1523075	20.000	ng/u1	0.00
88) Perylene-d12	24.069	264	1728932	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	41977	4.078	ng/uL	0.00
4) Pyridine-d5	3.793	84	570634	21.992	ng/u1	-0.01
7) Phenol-d5	7.211	99	767399	24.218	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	541057m	27.599	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.546	132	600450	25.103	ng/u1	-0.02
15) 4-Methylphenol-d8	8.764	113	638999	25.527	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	305056	25.677	ng/u1	0.00
24) 2-Nitrophenol-d4	9.946	143	342890	24.728	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	492368	19.926	ng/u1	0.02
31) 4-Chloroaniline-d4	11.040	131	725328	22.851	ng/u1	0.02
46) Dimethylphthalate-d6	14.087	166	1898343	26.118	ng/u1	0.00
49) Acenaphthylene-d8	14.363	160	2149381	26.306	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.010	143	182022m	18.977	ng/u1	0.05
60) Fluorene-d10	15.675	176	1585223	26.487	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	200649	17.105	ng/u1	0.01
73) Anthracene-d10	17.545	188	2307558	26.486	ng/u1	0.00
81) Pyrene-d10	19.769	212	2513717	25.275	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	2266542	25.988	ng/u1	-0.01

Target Compounds Qvalue

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

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