

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016127.D
 Acq On : 09 Jul 2023 11:42
 Operator : MA/JU
 Sample : 03356-06
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YBW35

Manual Integrations
 APPROVED

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

Quant Time: Jul 09 23:40:02 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	350999	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.846	136	1606192	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.669	164	1044639	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.433	188	2359194	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.539	240	2188380	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	2124918	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	39226	4.021	ng/uL	0.00
4) Pyridine-d5	3.793	84	501406	20.389	ng/u1	-0.01
7) Phenol-d5	7.216	99	656331	21.854	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	466513m	25.108	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.552	132	521856	23.019	ng/u1	-0.01
15) 4-Methylphenol-d8	8.769	113	549147	23.146	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	256186	20.870	ng/u1	0.00
24) 2-Nitrophenol-d4	9.951	143	296668	20.708	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	411848	16.132	ng/u1	0.04
31) 4-Chloroaniline-d4	11.046	131	640602	19.533	ng/u1	0.02
46) Dimethylphthalate-d6	14.087	166	1794087	22.220	ng/u1	0.00
49) Acenaphthylene-d8	14.363	160	1993502	21.963	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.028	143	185114m	17.373	ng/u1	0.06
60) Fluorene-d10	15.675	176	1548930	23.298	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	185003	12.751	ng/u1	0.02
73) Anthracene-d10	17.545	188	2463106	22.857	ng/u1	0.00
81) Pyrene-d10	19.769	212	2971264	20.793	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	2429115	22.662	ng/u1	-0.01

Target Compounds Qvalue

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

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