

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011092.D
 Acq On : 25 Jul 2022 15:01
 Operator : CG/JU
 Sample : N3801-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHX8

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/26/2022
 Supervised By :mohammad ahmed 07/26/2022

Quant Time: Jul 25 23:21:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	430066	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.463	136	1991978	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.334	164	1223594	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	2508127	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.216	240	2104499	20.000	ng/u1	#-0.01
88) Perylene-d12	23.563	264	1791502	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	52158	4.372	ng/uL	-0.02
4) Pyridine-d5	3.575	84	338730	10.985	ng/u1	-0.02
7) Phenol-d5	6.846	99	251345	7.120	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.010	67	912691	40.205	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.199	132	818736	28.702	ng/u1	-0.01
15) 4-Methylphenol-d8	8.381	113	530657	18.776	ng/u1	-0.02
21) Nitrobenzene-d5	8.828	128	587751	40.488	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.551	143	585471	40.759	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	885735	33.718	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.610	131	1422394	32.845	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	3637365	43.494	ng/u1	-0.01
49) Acenaphthylene-d8	14.022	160	4220658	42.775	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	76273m	4.506	ng/u1	0.00
60) Fluorene-d10	15.334	176	3004859	42.272	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	497160	40.058	ng/u1	0.00
73) Anthracene-d10	17.198	188	4666017	43.760	ng/u1	0.00
81) Pyrene-d10	19.451	212	5188982	46.947	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.415	264	3794588	44.100	ng/u1	-0.01

Target Compounds Qvalue

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

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