

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011086.D
 Acq On : 25 Jul 2022 11:34
 Operator : CG/JU
 Sample : N3801-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHX2

Manual Integrations
 APPROVED

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

Quant Time: Jul 25 23:19:58 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	7.669	152	449654	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.463	136	1987167	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.334	164	1111435	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	1956630	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.216	240	1453272	20.000	ng/u1	#-0.01
88) Perylene-d12	23.563	264	1542973	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	60745	4.870	ng/uL	-0.02
4) Pyridine-d5	3.581	84	214684	6.659	ng/u1	-0.01
7) Phenol-d5	6.846	99	233636	6.330	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	833811	35.131	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.199	132	747476	25.062	ng/u1	-0.01
15) 4-Methylphenol-d8	8.381	113	477006	16.143	ng/u1	-0.02
21) Nitrobenzene-d5	8.834	128	520727	35.958	ng/u1	0.00
24) 2-Nitrophenol-d4	9.552	143	412267	28.771	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	745935	28.465	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.610	131	970336	22.460	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	2715938	35.754	ng/u1	-0.01
49) Acenaphthylene-d8	14.022	160	3405296	37.994	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	34979m	2.275	ng/u1	0.00
60) Fluorene-d10	15.334	176	2239432	34.683	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	217525	22.467	ng/u1	0.00
73) Anthracene-d10	17.198	188	3070039	36.908	ng/u1	0.00
81) Pyrene-d10	19.451	212	3044840	39.893	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.410	264	2725218	36.774	ng/u1	-0.02

Target Compounds Qvalue

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

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