

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009892.D
 Acq On : 16 Apr 2022 09:34
 Operator : CG/JU
 Sample : N2421-10
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY7

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022

Quant Time: Apr 18 01:10:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	138937	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	622670	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	442694	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.339	188	960951	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.416	240	982665	20.000	ng/u1	# 0.00
88) Perylene-d12	23.880	264	755973	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	15126m	4.859	ng/uL	0.00
4) Pyridine-d5	3.799	84	93749	10.922	ng/u1	0.00
7) Phenol-d5	7.111	99	108407	8.887	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	268738	37.007	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	272545	29.621	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	198963	19.540	ng/u1	0.00
21) Nitrobenzene-d5	9.110	128	178670	39.791	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	194019	39.412	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	366461m	35.574	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	294595	20.078	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	1409308	40.527	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1533515	37.223	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	55004	8.722	ng/u1	0.00
60) Fluorene-d10	15.581	176	1212180	41.437	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	216815	38.086	ng/u1	0.00
73) Anthracene-d10	17.439	188	1980669	43.049	ng/u1	0.00
81) Pyrene-d10	19.669	212	2436155	45.393	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	1857840	47.191	ng/u1	0.00
Target Compounds						
8) Phenol	7.134	94	19265	1.578	ng/u1#	85
86) Bis(2-ethylhexyl)phtha...	21.327	149	49898	1.291	ng/u1#	83

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

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