

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010858.D
 Acq On : 26 Jun 2022 09:04
 Operator : CG/JU
 Sample : N3392-20 10X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEZ5

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022

Quant Time: Jun 26 22:57:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	475694	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1920049	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1030199	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1989276	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1978589	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	2051902	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	34465	2.741	ng/uL	0.00
4) Pyridine-d5	3.616	84	229215m	6.848	ng/u1	0.01
7) Phenol-d5	6.893	99	278775	7.403	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	907756	36.860	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	911590	29.178	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	532483	17.888	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	547730	44.812	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	497748	49.222	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	848391	33.954	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.651	131	1221222	29.942	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2621819	37.152	ng/u1	-0.01
49) Acenaphthylene-d8	14.063	160	3290899	36.539	ng/u1	0.00
54) 4-Nitrophenol-d4	14.581	143	68822	6.244	ng/u1	0.00
60) Fluorene-d10	15.369	176	2301658	38.007	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	191392	29.040	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3626461	40.974	ng/u1	-0.01
81) Pyrene-d10	19.480	212	4098306	38.020	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.462	264	4287868	42.288	ng/u1	-0.01

Target Compounds Qvalue

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

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