

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010856.D
 Acq On : 26 Jun 2022 07:56
 Operator : CG/JU
 Sample : N3392-11 10X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY6

Quant Time: Jun 26 22:57:38 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	7.710	152	474060	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	1900159	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	988922	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1758362	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1695932	20.000	ng/u1	#-0.01
88) Perylene-d12	23.609	264	1815675	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	35034	2.796	ng/uL	0.00
4) Pyridine-d5	3.616	84	294806	8.837	ng/u1	0.01
7) Phenol-d5	6.893	99	305867	8.151	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	948710	38.656	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	950331	30.523	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	572000	19.282	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	571872	47.276	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	528596	52.819	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.134	165	869551	35.165	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1301271	32.239	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2879785	42.511	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3489587	40.362	ng/u1	0.00
54) 4-Nitrophenol-d4	14.581	143	71523	6.760	ng/u1	0.00
60) Fluorene-d10	15.369	176	2415450	41.550	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	193545	33.223	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3517238	44.959	ng/u1	-0.01
81) Pyrene-d10	19.486	212	3879841	41.992	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.462	264	4240332	47.260	ng/u1	-0.01

Target Compounds Qvalue

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

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