

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010860.D
 Acq On : 26 Jun 2022 10:11
 Operator : CG/JU
 Sample : N3392-12 10X
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY7

Quant Time: Jun 26 22:58:13 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	471114	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	1911903	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1008309	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1880776	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1858498	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1927827	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	52749	4.235	ng/uL	0.00
4) Pyridine-d5	3.611	84	654804	19.752	ng/u1	0.00
7) Phenol-d5	6.887	99	588943	15.792	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	1028402	42.165	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	1082978	35.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	747931	25.370	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	620816	51.007	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	570413	56.648	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	973468	39.126	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.651	131	1403946	34.569	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	3184186	46.101	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	3821393	43.350	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.581	143	157178	14.569	ng/u1	0.00
60) Fluorene-d10	15.369	176	2688197	45.353	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	231853	37.208	ng/u1	-0.01
73) Anthracene-d10	17.233	188	4017506	48.011	ng/u1	-0.01
81) Pyrene-d10	19.480	212	4475804	44.205	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	4810339	50.494	ng/u1	-0.02

Target Compounds Qvalue

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

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