

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010861.D
 Acq On : 26 Jun 2022 10:45
 Operator : CG/JU
 Sample : N3392-18 10X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEZ3

Quant Time: Jun 26 22:58:22 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	527380	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	2152527	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1129390	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	2031853	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1845072	20.000	ng/u1	#-0.01
88) Perylene-d12	23.609	264	1986203	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	36663	2.630	ng/uL	0.00
4) Pyridine-d5	3.616	84	296628	7.993	ng/u1	0.01
7) Phenol-d5	6.887	99	343273	8.223	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	1041113	38.132	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	1045407	30.182	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	641770	19.446	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	628033	45.832	ng/u1	0.00
24) 2-Nitrophenol-d4	9.592	143	577671	50.956	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.134	165	960714	34.297	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1299953	28.430	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	3046939	39.385	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	3743059	37.909	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.580	143	76002	6.290	ng/u1	0.00
60) Fluorene-d10	15.369	176	2573029	38.756	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	206179	30.628	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3726562	41.223	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3955708	39.352	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.456	264	4191784	42.708	ng/u1	-0.02

Target Compounds Qvalue

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

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