

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010864.D
 Acq On : 26 Jun 2022 12:26
 Operator : CG/JU
 Sample : N3392-16 10X
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEZ1

Quant Time: Jun 26 22:58:47 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	558524	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.504	136	2294522	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.369	164	1240010	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.134	188	2170165	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1841238	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	2009220	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	30383	2.058	ng/uL	0.00
4) Pyridine-d5	3.617	84	310012	7.888	ng/u1	0.01
7) Phenol-d5	6.887	99	314041	7.103	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	834625	28.865	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	850629	23.189	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	545145	15.597	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	501251	34.316	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	448258	37.093	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	770870	25.817	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.652	131	1174260	24.092	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2497057	29.397	ng/u1	-0.01
49) Acenaphthylene-d8	14.057	160	3041847	28.059	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.581	143	69123	5.210	ng/u1	0.00
60) Fluorene-d10	15.369	176	2163076	29.675	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.487	200	157657	21.927	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3168166	32.812	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3233451	32.234	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3342912	33.669	ng/u1	-0.02

Target Compounds Qvalue

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

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