

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010862.D
 Acq On : 26 Jun 2022 11:19
 Operator : CG/JU
 Sample : N3392-08 10X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY3

Quant Time: Jun 26 22:58:30 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	522221	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	2111738	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	1130694	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	2175612	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	2176183	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	2309314	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	31079	2.251	ng/uL	0.00
4) Pyridine-d5	3.616	84	244851	6.663	ng/u1	0.01
7) Phenol-d5	6.887	99	263435	6.372	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	821604	30.390	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	822003	23.967	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	495497	15.162	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	482550	35.895	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	441752	39.719	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	766857	27.905	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.651	131	1107156	24.681	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2447032	31.594	ng/u1	-0.01
49) Acenaphthylene-d8	14.057	160	3018427	30.535	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.581	143	65631	5.425	ng/u1	0.00
60) Fluorene-d10	15.369	176	2105936	31.684	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	168530	23.381	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3223303	33.300	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3663027	30.896	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3855557	33.786	ng/u1	-0.02

Target Compounds Qvalue

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

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