

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010847.D
 Acq On : 26 Jun 2022 02:19
 Operator : CG/JU
 Sample : N3392-19 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEZ4

Quant Time: Jun 26 22:55:52 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	7.716	152	465905	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1892705	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	992607	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1786387	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1625707	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1779259	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	36060	2.928	ng/uL	0.00
4) Pyridine-d5	3.617	84	421785	12.865	ng/u1	0.01
7) Phenol-d5	6.893	99	331055	8.976	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.058	67	972694	40.327	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	974203	31.838	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	603122	20.687	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	590238	48.987	ng/u1	0.00
24) 2-Nitrophenol-d4	9.599	143	540594	54.231	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	909019	36.906	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1445917	35.963	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2947614	43.351	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	3591798	41.390	ng/u1	0.00
54) 4-Nitrophenol-d4	14.581	143	73449	6.916	ng/u1	0.00
60) Fluorene-d10	15.369	176	2486384	42.612	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	193300	32.660	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3515593	44.233	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3724724	42.054	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3978201	45.246	ng/u1	-0.02

Target Compounds Qvalue

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

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