

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010876.D
 Acq On : 27 Jun 2022 14:16
 Operator : CG/JU
 Sample : N3392-20 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jun 27 23:15:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 27 23:12:49 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	411627	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	1633499	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	822995	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	1476186	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1372266	20.000	ng/u1	# 0.00
88) Perylene-d12	23.604	264	1546509	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	32558	2.992	ng/uL	0.00
4) Pyridine-d5	3.617	84	207443	7.162	ng/u1	0.00
7) Phenol-d5	6.887	99	258453	7.932	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	860178	40.365	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	851256	31.488	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	495209	19.225	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	497235	47.817	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	450340	52.346	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	770578	36.250	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	1106643	31.892	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2207909	39.164	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	2833533	39.381	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	51424	5.840	ng/u1	0.00
60) Fluorene-d10	15.363	176	1945056	40.204	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.487	200	179538	36.710	ng/u1	0.00
73) Anthracene-d10	17.228	188	2914372	44.374	ng/u1	0.00
81) Pyrene-d10	19.480	212	3157819	42.239	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	3476387	45.489	ng/u1	0.00

Target Compounds Qvalue

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

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