

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010878.D
 Acq On : 27 Jun 2022 15:24
 Operator : CG/JU
 Sample : N3392-12 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jun 27 23:15:48 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	7.711	152	470386	20.000	ng/u1	0.00
20) Naphthalene-d8	10.504	136	1925962	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.363	164	979652	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	1673083	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1366932	20.000	ng/u1	# 0.00
88) Perylene-d12	23.604	264	1521826	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	52485	4.221	ng/uL	0.00
4) Pyridine-d5	3.611	84	653545	19.744	ng/u1	0.00
7) Phenol-d5	6.887	99	575454	15.454	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	1011707	41.545	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	1042945	33.759	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	735177	24.976	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	599071	48.861	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	548978	54.121	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	936891	37.381	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	1376948	33.656	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2965861	44.196	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	3608802	42.136	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	128414	12.251	ng/u1	0.00
60) Fluorene-d10	15.363	176	2512189	43.623	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	232887	42.014	ng/u1	0.00
73) Anthracene-d10	17.228	188	3500878	47.031	ng/u1	0.00
81) Pyrene-d10	19.480	212	3562007	47.831	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	3662204	48.698	ng/u1	0.00

Target Compounds Qvalue

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

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