

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010859.D
 Acq On : 26 Jun 2022 09:38
 Operator : CG/JU
 Sample : N3392-02 10X
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEX7

Quant Time: Jun 26 22:58:04 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	439810	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	1752837	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	926929	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1757683	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1701682	20.000	ng/u1	#-0.01
88) Perylene-d12	23.610	264	1770471	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	27478	2.363	ng/uL	0.00
4) Pyridine-d5	3.617	84	309980	10.016	ng/u1	0.01
7) Phenol-d5	6.887	99	236893	6.804	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	770106	33.822	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	759006	26.277	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	448774	16.306	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	456589	40.918	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	406534	44.037	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.134	165	711931	31.211	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1034592	27.786	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2397801	37.764	ng/u1	-0.01
49) Acenaphthylene-d8	14.057	160	2877078	35.503	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.581	143	60196	6.070	ng/u1	0.00
60) Fluorene-d10	15.369	176	2110943	38.741	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	174978	30.048	ng/u1	-0.01
73) Anthracene-d10	17.233	188	3255677	41.632	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3645852	39.326	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3783516	43.246	ng/u1	-0.02

Target Compounds Qvalue

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

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