

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071122\
 Data File : BP010958.D
 Acq On : 11 Jul 2022 20:25
 Operator : CG/JU
 Sample : N3625-02 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0ZB9

Quant Time: Jul 12 01:06:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 06 23:13:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	520529	20.000	ng/u1	0.00
20) Naphthalene-d8	10.493	136	2111088	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.351	164	1040621	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	1764002	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1497863	20.000	ng/u1	#-0.01
88) Perylene-d12	23.586	264	1710120	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	60156	4.372	ng/uL	0.00
4) Pyridine-d5	3.611	84	311231	8.497	ng/u1	0.00
7) Phenol-d5	6.875	99	286377	6.950	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	798075	29.615	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	793073	23.198	ng/u1	0.00
15) 4-Methylphenol-d8	8.411	113	514398	15.792	ng/u1	-0.01
21) Nitrobenzene-d5	8.858	128	456612	33.976	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	412411	37.092	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	714729	26.016	ng/u1	0.00
31) 4-Chloroaniline-d4	10.634	131	911718	20.331	ng/u1	0.00
46) Dimethylphthalate-d6	13.775	166	2319564	32.540	ng/u1	0.00
49) Acenaphthylene-d8	14.046	160	2786983	30.634	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	56326	5.059	ng/u1	0.00
60) Fluorene-d10	15.351	176	1958671	32.019	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	184704	31.604	ng/u1	0.00
73) Anthracene-d10	17.216	188	2751720	35.061	ng/u1	0.00
81) Pyrene-d10	19.463	212	2839929	34.801	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.433	264	3140382	37.161	ng/u1	-0.01

Target Compounds	Qvalue

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

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