

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071122\
 Data File : BP010948.D
 Acq On : 11 Jul 2022 13:09
 Operator : CG/JU
 Sample : N3639-01 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0L62

Quant Time: Jul 11 13:41:56 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	7.698	152	227771	20.000	ng/u1	0.00
20) Naphthalene-d8	10.492	136	923406	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.357	164	459321	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	794718	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.233	240	673594	20.000	ng/u1	# 0.00
88) Perylene-d12	23.592	264	751741	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	38423	6.381	ng/uL	0.00
4) Pyridine-d5	3.610	84	447693	27.932	ng/u1	0.00
7) Phenol-d5	6.875	99	341356	18.932	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	915244	77.617	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	921853	61.624	ng/u1	0.00
15) 4-Methylphenol-d8	8.410	113	592428	41.564	ng/u1	-0.01
21) Nitrobenzene-d5	8.863	128	520878	88.609	ng/u1	0.00
24) 2-Nitrophenol-d4	9.581	143	466862	95.997	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	806122	67.084	ng/u1	0.00
31) 4-Chloroaniline-d4	10.639	131	459510	23.426	ng/u1	0.00
46) Dimethylphthalate-d6	13.774	166	2472714	78.589	ng/u1	0.00
49) Acenaphthylene-d8	14.045	160	3085711	76.842	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	63357	12.892	ng/u1	0.00
60) Fluorene-d10	15.357	176	2147490	79.534	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.474	200	201857	76.665	ng/u1	0.00
73) Anthracene-d10	17.221	188	3149140	89.064	ng/u1	0.00
81) Pyrene-d10	19.468	212	3242104	88.346	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.445	264	3559638	95.824	ng/u1	0.00

Target Compounds Qvalue

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

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